



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 01:30 AM UTC

PDB ID : 5VYE / pdb_00005vye
Title : Crystal Structure of L-Threonine Aldolase from Pseudomonas putida
Authors : Beaudoin, S.F.; Burg, M.J.; Stewart, J.D.
Deposited on : 2017-05-25
Resolution : 2.27 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

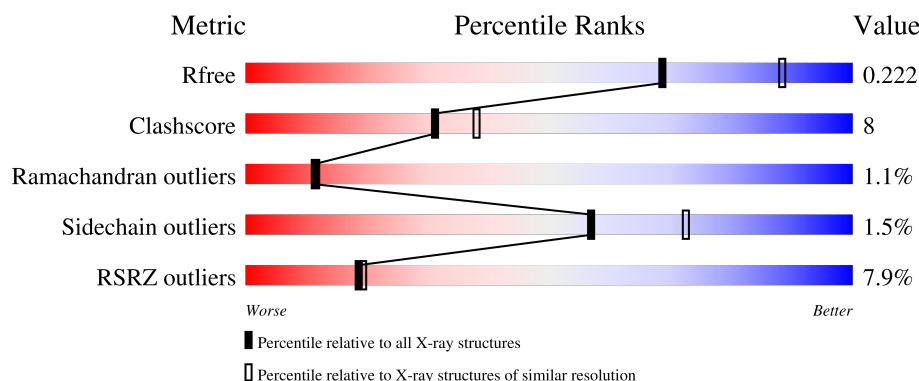
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.27 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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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	346	8% 87% 11% ..
1	B	346	7% 85% 14% ..
1	C	346	7% 82% 16% ..
1	D	346	9% 85% 12% ..

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
2	PLR	A	401	-	X	-	-
2	PLR	B	401	-	X	-	-
2	PLR	C	401	-	-	X	-
2	PLR	D	401	-	X	X	-

2 Entry composition [i](#)

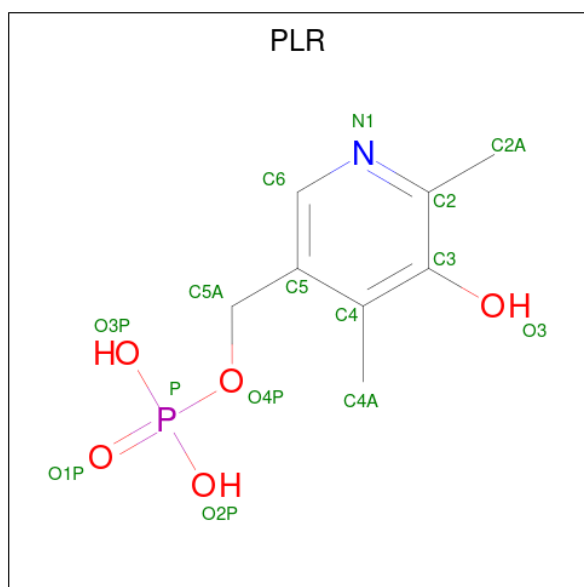
There are 4 unique types of molecules in this entry. The entry contains 21529 atoms, of which 10299 are hydrogens and 0 are deuteriums.

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 L-threonine aldolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	343	Total	C	H	N	O	S	0	4	0
			5239	1683	2566	467	504	19			
1	B	344	Total	C	H	N	O	S	0	4	0
			5251	1688	2569	466	509	19			
1	C	342	Total	C	H	N	O	S	0	2	0
			5202	1674	2546	462	501	19			
1	D	342	Total	C	H	N	O	S	0	2	0
			5202	1674	2546	462	501	19			

- Molecule 2 is (5-HYDROXY-4,6-DIMETHYLPYRIDIN-3-YL)METHYL DIHYDROGEN PHOSPHATE (CCD ID: PLR) (formula: $C_8H_{12}NO_5P$).



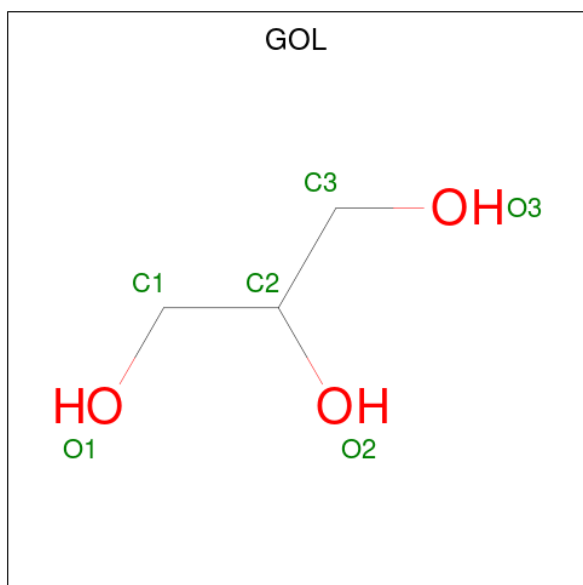
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	0	0
			25	8	10	1	5	1		
2	B	1	Total	C	H	N	O	P	0	0
			25	8	10	1	5	1		

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	C	1	Total	C	H	N	O	P	0	0
			25	8	10	1	5	1		
2	D	1	Total	C	H	N	O	P	0	0
			25	8	10	1	5	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	B	1	Total	C	H	O	0	0
			14	3	8	3		
3	C	1	Total	C	H	O	0	0
			14	3	8	3		
3	D	1	Total	C	H	O	0	0
			14	3	8	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	134	Total	O	0	0
			134	134		
4	B	109	Total	O	0	0
			109	109		
4	C	124	Total	O	0	0
			124	124		

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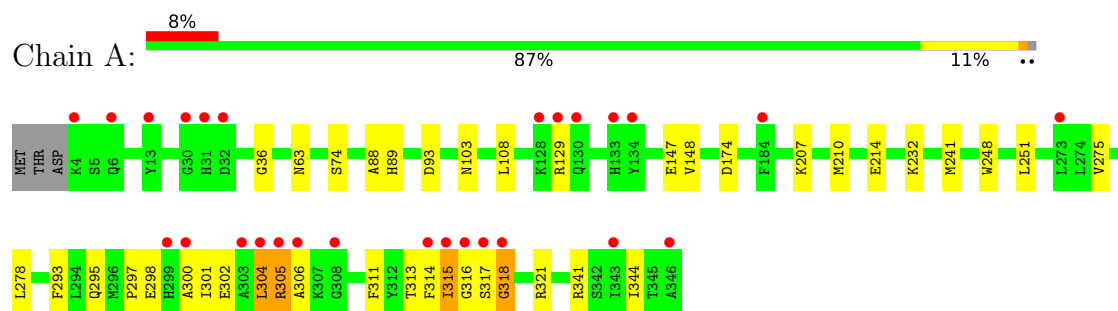
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	112	Total 112	O 112	0	0

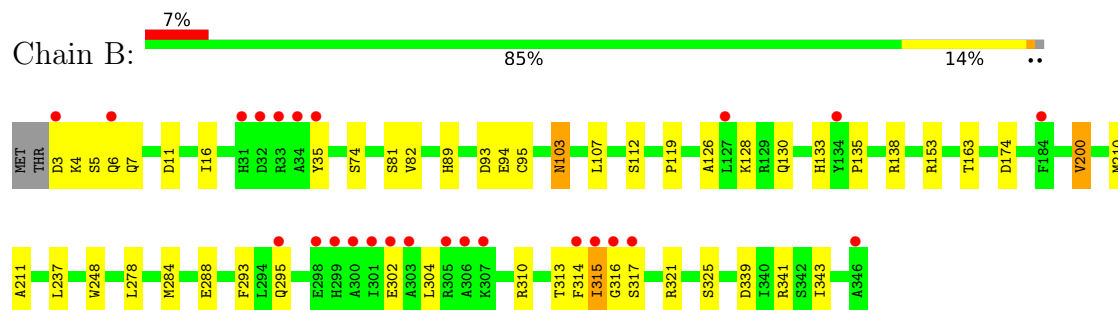
3 Residue-property plots [i](#)

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.

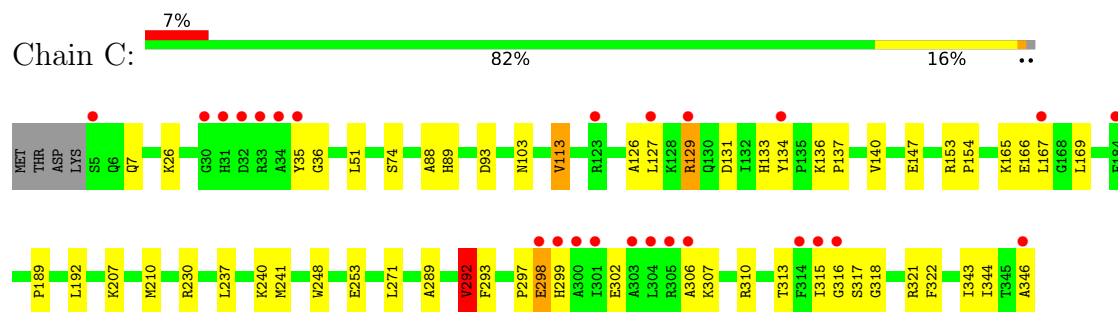
• Molecule 1: L-threonine aldolase



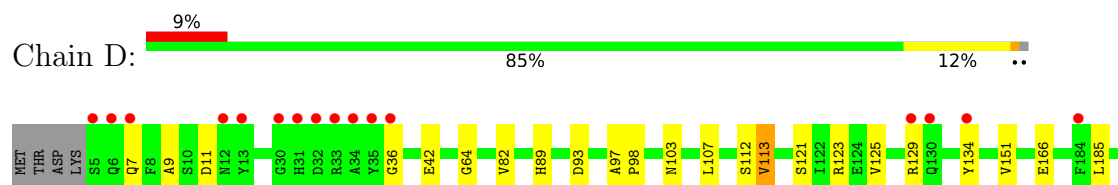
• Molecule 1: L-threonine aldolase

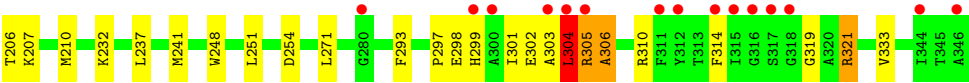


• Molecule 1: L-threonine aldolase



• Molecule 1: L-threonine aldolase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	198.17Å 186.96Å 53.27Å 90.00° 98.92° 90.00°	Depositor
Resolution (Å)	26.75 – 2.27 26.75 – 2.27	Depositor EDS
% Data completeness (in resolution range)	99.3 (26.75-2.27) 94.4 (26.75-2.27)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.29Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.180 , 0.219 0.182 , 0.222	Depositor DCC
R_{free} test set	2000 reflections (2.28%)	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	21529	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.64	0/2749	0.74	0/3724
1	B	0.68	1/2753 (0.0%)	0.72	0/3729
1	C	0.65	1/2723 (0.0%)	0.75	2/3690 (0.1%)
1	D	0.63	1/2723 (0.0%)	0.71	0/3690
All	All	0.65	3/10948 (0.0%)	0.73	2/14833 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	200	VAL	CB-CG1	-7.08	1.29	1.52
1	D	11	ASP	CA-C	6.55	1.61	1.52
1	C	292	VAL	CB-CG1	-5.45	1.34	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	113	VAL	CA-CB-CG2	5.68	120.06	110.40
1	C	113	VAL	CG1-CB-CG2	5.24	122.33	110.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2673	2566	2548	37	2
1	B	2682	2569	2559	50	0
1	C	2656	2546	2538	52	2
1	D	2656	2546	2538	41	0
2	A	15	10	10	3	0
2	B	15	10	10	4	0
2	C	15	10	10	8	0
2	D	15	10	10	6	0
3	A	6	8	8	1	0
3	B	6	8	8	1	0
3	C	6	8	8	2	0
3	D	6	8	8	0	0
4	A	134	0	0	0	0
4	B	109	0	0	8	0
4	C	124	0	0	10	0
4	D	112	0	0	4	0
All	All	11230	10299	10255	173	2

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 8.

All (173) 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:301:ILE:O	1:A:305:ARG:HB2	1.78	0.83
1:A:301:ILE:HG22	1:A:305:ARG:HD2	1.66	0.77
1:B:103:ASN:OD1	1:D:103:ASN:HB2	1.84	0.77
1:B:103:ASN:HB2	1:D:103:ASN:ND2	2.01	0.75
1:B:153:ARG:NH2	1:B:288:GLU:OE2	2.21	0.74
1:B:126:ALA:O	1:B:130:GLN:NE2	2.24	0.69
2:C:401:PLR:H4A3	2:C:401:PLR:O4P	1.96	0.66
1:D:304:LEU:O	1:D:306:ALA:N	2.29	0.65
1:A:302:GLU:HA	1:A:305:ARG:HB2	1.79	0.64
1:C:297:PRO:O	1:C:299:HIS:N	2.34	0.61
1:B:304:LEU:HD22	4:B:609:HOH:O	2.01	0.61
1:C:237:LEU:HD13	4:C:509:HOH:O	2.00	0.60
1:B:89:HIS:CE1	2:B:401:PLR:H5A1	2.37	0.60
1:A:103:ASN:ND2	1:C:103:ASN:HB2	2.16	0.60
1:D:297:PRO:O	1:D:299:HIS:N	2.35	0.60
1:B:302:GLU:HG2	1:B:343:ILE:HG23	1.85	0.59
2:D:401:PLR:O4P	2:D:401:PLR:H4A3	2.02	0.59
1:B:284:MET:HG2	1:B:295:GLN:HG3	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:297:PRO:C	1:C:299:HIS:H	2.12	0.57
1:D:123:ARG:NH2	1:D:166:GLU:OE1	2.38	0.57
1:A:293:PHE:CZ	1:A:321:ARG:HD2	2.40	0.57
1:A:210:MET:HE3	1:A:251:LEU:HD22	1.87	0.56
1:C:271:LEU:HD23	1:C:292:VAL:HG11	1.86	0.56
1:A:298:GLU:C	1:A:300:ALA:H	2.13	0.56
1:D:7:GLN:HG2	1:D:9:ALA:H	1.71	0.55
1:B:5:SER:HB2	4:B:588:HOH:O	2.06	0.55
1:C:306:ALA:O	1:C:307:LYS:HB3	2.06	0.55
1:A:298:GLU:C	1:A:300:ALA:N	2.62	0.55
1:B:314:PHE:C	1:B:316:GLY:H	2.14	0.55
1:C:298:GLU:O	1:C:302:GLU:HG2	2.07	0.54
1:D:89:HIS:CD2	1:D:93:ASP:HB2	2.42	0.54
1:C:127:LEU:N	1:C:127:LEU:HD12	2.23	0.54
1:C:26:LYS:NZ	4:C:507:HOH:O	2.39	0.54
1:D:314:PHE:N	1:D:319:GLY:O	2.41	0.53
1:C:316:GLY:O	1:C:318:GLY:N	2.40	0.53
1:D:113:VAL:HG23	4:D:556:HOH:O	2.09	0.53
1:B:315:ILE:CG2	4:C:615:HOH:O	2.56	0.53
1:A:304:LEU:HG	1:A:311:PHE:CZ	2.44	0.53
1:B:313:THR:HG21	4:B:609:HOH:O	2.08	0.53
1:C:293:PHE:CZ	1:C:321:ARG:HD2	2.43	0.53
1:D:304:LEU:HD23	1:D:305:ARG:HG3	1.91	0.53
1:C:133:HIS:CE1	4:C:553:HOH:O	2.62	0.52
1:B:138:ARG:CG	4:B:602:HOH:O	2.58	0.52
1:A:301:ILE:O	1:A:305:ARG:N	2.39	0.52
1:A:210:MET:HE1	1:A:248:TRP:CD2	2.44	0.52
1:C:129:ARG:HD3	1:C:134:TYR:CD2	2.45	0.51
1:C:207:LYS:HZ3	2:C:401:PLR:C4A	2.23	0.51
1:D:129:ARG:O	1:D:129:ARG:HG3	2.10	0.51
1:C:210:MET:HE1	1:C:248:TRP:CD2	2.46	0.51
1:B:89:HIS:CD2	1:B:93:ASP:HB2	2.46	0.51
1:B:313:THR:CG2	4:B:609:HOH:O	2.59	0.51
1:C:292:VAL:HG13	1:C:322:PHE:HB2	1.93	0.51
1:D:129:ARG:NH1	1:D:134:TYR:CD1	2.79	0.51
1:B:133:HIS:CE1	4:B:594:HOH:O	2.63	0.51
1:B:35:TYR:CD1	1:B:237:LEU:HD23	2.47	0.50
1:A:103:ASN:HB2	1:C:103:ASN:OD1	2.13	0.49
1:D:36:GLY:HA2	1:D:241:MET:HG2	1.93	0.49
1:B:138:ARG:HG2	4:B:602:HOH:O	2.11	0.49
1:C:89:HIS:CD2	1:C:93:ASP:HB2	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:MET:HE1	1:B:248:TRP:CD2	2.48	0.49
1:C:237:LEU:CD1	4:C:509:HOH:O	2.58	0.49
1:B:3:ASP:CB	1:B:4:LYS:HB2	2.43	0.49
1:D:232:LYS:HD3	1:D:237:LEU:HD22	1.94	0.49
1:D:302:GLU:O	1:D:306:ALA:HB3	2.13	0.48
1:B:278:LEU:HD11	1:B:341:ARG:CZ	2.43	0.48
1:B:4:LYS:HG3	1:B:6:GLN:H	1.79	0.48
1:A:36:GLY:HA2	1:A:241:MET:HG2	1.96	0.48
1:D:232:LYS:NZ	4:D:508:HOH:O	2.46	0.48
1:A:301:ILE:HG22	1:A:305:ARG:CD	2.41	0.47
1:A:316:GLY:HA2	1:A:317:SER:HA	1.70	0.47
1:B:7:GLN:HB3	1:B:310:ARG:HB2	1.97	0.47
1:B:302:GLU:HG2	1:B:343:ILE:CG2	2.44	0.47
1:A:89:HIS:CE1	2:A:401:PLR:C5	2.97	0.47
1:D:304:LEU:HD23	1:D:305:ARG:H	1.80	0.47
1:D:210:MET:HE1	1:D:248:TRP:CD2	2.50	0.47
1:B:315:ILE:HG22	4:C:615:HOH:O	2.15	0.46
1:D:303:ALA:O	1:D:304:LEU:O	2.33	0.46
1:B:103:ASN:OD1	1:D:103:ASN:CB	2.58	0.46
1:D:207:LYS:NZ	2:D:401:PLR:H4A1	2.30	0.46
1:A:174:ASP:OD2	2:A:401:PLR:N1	2.49	0.46
1:A:297:PRO:O	1:A:300:ALA:N	2.43	0.46
1:C:89:HIS:CE1	2:C:401:PLR:H5A1	2.50	0.46
1:C:89:HIS:CD2	2:C:401:PLR:C3	2.98	0.46
1:C:297:PRO:HD2	1:C:299:HIS:ND1	2.31	0.45
1:C:298:GLU:O	1:C:298:GLU:HG3	2.16	0.45
1:A:304:LEU:HD11	1:A:313:THR:CG2	2.47	0.45
1:B:7:GLN:CG	1:B:310:ARG:HB2	2.45	0.45
1:B:11:ASP:HA	1:B:325:SER:OG	2.16	0.45
1:A:278:LEU:HD21	1:A:341:ARG:NH1	2.32	0.45
1:C:129:ARG:HB2	1:C:134:TYR:CD1	2.51	0.45
1:C:240:LYS:HE3	1:D:206:THR:HG22	1.98	0.45
1:C:292:VAL:CG1	1:C:322:PHE:HB2	2.46	0.45
1:A:232:LYS:HG2	1:B:95:CYS:HB2	1.99	0.45
1:A:108:LEU:HD13	1:A:129:ARG:HD3	1.98	0.45
1:A:298:GLU:O	1:A:301:ILE:HG13	2.17	0.45
1:B:293:PHE:CZ	1:B:321:ARG:HD2	2.51	0.45
1:C:230:ARG:HH21	3:C:402:GOL:C3	2.29	0.45
1:C:88:ALA:HA	1:C:147:GLU:OE2	2.16	0.45
1:A:210:MET:HE1	1:A:248:TRP:CE2	2.52	0.45
1:B:210:MET:HE1	1:B:248:TRP:CE2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:129:ARG:O	1:C:129:ARG:HG3	2.18	0.44
1:B:4:LYS:HD2	4:B:588:HOH:O	2.16	0.44
1:B:74:SER:O	3:B:402:GOL:H32	2.17	0.44
1:C:127:LEU:HD11	1:C:167:LEU:CD2	2.48	0.44
1:B:313:THR:O	1:B:314:PHE:HB2	2.18	0.44
1:B:3:ASP:HB3	1:B:4:LYS:HB2	2.00	0.44
1:B:119:PRO:HB3	1:B:163:THR:OG1	2.18	0.44
1:B:174:ASP:OD2	2:B:401:PLR:N1	2.51	0.44
1:A:295:GLN:OE1	1:A:318:GLY:HA2	2.18	0.44
1:D:123:ARG:NH2	1:D:166:GLU:CD	2.76	0.44
1:A:89:HIS:CD2	1:A:93:ASP:HB2	2.53	0.43
1:C:126:ALA:HA	1:C:137:PRO:HG2	2.00	0.43
1:C:7:GLN:HG3	1:C:310:ARG:HB2	2.01	0.43
1:A:63:ASN:C	1:A:214[B]:GLU:HG3	2.43	0.43
1:A:300:ALA:C	1:A:302:GLU:N	2.75	0.43
1:D:207:LYS:HZ3	2:D:401:PLR:C4A	2.31	0.43
1:A:210:MET:CE	1:A:251:LEU:HD22	2.49	0.43
1:B:82:VAL:HB	1:B:107:LEU:HD23	2.00	0.43
1:A:148:VAL:HA	1:A:293:PHE:CZ	2.54	0.43
1:B:93:ASP:HA	1:C:133:HIS:HB3	2.00	0.43
1:C:35:TYR:HE1	4:C:545:HOH:O	2.02	0.43
1:B:314:PHE:CB	1:B:316:GLY:H	2.32	0.43
1:C:140:VAL:HG23	1:C:169:LEU:HD13	2.01	0.43
1:A:88:ALA:HA	1:A:147:GLU:OE2	2.19	0.42
1:C:127:LEU:C	1:C:129:ARG:H	2.27	0.42
1:D:304:LEU:O	1:D:305:ARG:C	2.62	0.42
1:D:113:VAL:O	1:D:113:VAL:CG2	2.68	0.42
1:C:51:LEU:HG	1:C:189:PRO:HG2	2.02	0.42
1:C:127:LEU:N	1:C:127:LEU:CD1	2.82	0.42
1:D:97:ALA:HB3	1:D:98:PRO:HD3	2.02	0.42
1:D:151:VAL:CG1	1:D:185:LEU:HD11	2.49	0.42
1:C:299:HIS:HB3	1:C:343:ILE:CG2	2.50	0.42
1:C:89:HIS:CE1	2:C:401:PLR:C5	3.03	0.42
1:D:254:ASP:O	1:D:254:ASP:CG	2.63	0.42
1:D:310:ARG:HD3	1:D:310:ARG:HA	1.86	0.42
1:D:82:VAL:HB	1:D:107:LEU:HD23	2.02	0.42
1:A:305:ARG:HB3	1:A:306:ALA:H	1.54	0.41
1:A:317:SER:O	1:A:318:GLY:C	2.63	0.41
1:D:121:SER:O	4:D:501:HOH:O	2.21	0.41
1:D:210:MET:HE3	1:D:251:LEU:HD22	2.01	0.41
1:D:293:PHE:CZ	1:D:321:ARG:HD2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:LYS:NZ	2:A:401:PLR:O3	2.53	0.41
1:B:339:ASP:O	1:B:343:ILE:HD12	2.21	0.41
1:B:16:ILE:HD11	1:B:211:ALA:HB2	2.01	0.41
1:C:74:SER:HA	3:C:402:GOL:H31	2.03	0.41
1:B:89:HIS:CE1	2:B:401:PLR:C5A	3.03	0.41
1:B:89:HIS:CE1	1:B:94:GLU:HG2	2.56	0.41
1:C:136:LYS:CE	4:C:510:HOH:O	2.68	0.41
1:B:7:GLN:HG3	1:B:310:ARG:HB2	2.03	0.41
1:C:165:LYS:O	1:C:166:GLU:C	2.62	0.41
1:C:253:GLU:OE2	4:C:501:HOH:O	2.22	0.41
1:D:237:LEU:HD13	4:D:513:HOH:O	2.20	0.41
1:D:297:PRO:C	1:D:299:HIS:N	2.79	0.41
1:A:74:SER:O	3:A:402:GOL:H32	2.21	0.41
1:A:298:GLU:O	1:A:300:ALA:N	2.54	0.41
1:C:36:GLY:HA2	1:C:241:MET:HG2	2.02	0.41
1:C:131:ASP:HB2	4:C:597:HOH:O	2.20	0.41
1:C:153:ARG:O	1:C:154:PRO:C	2.64	0.41
1:C:207:LYS:NZ	2:C:401:PLR:C4A	2.84	0.41
1:B:89:HIS:CE1	2:B:401:PLR:C5	3.04	0.40
1:C:207:LYS:HZ3	2:C:401:PLR:C4	2.34	0.40
1:C:207:LYS:NZ	2:C:401:PLR:H4A1	2.36	0.40
1:A:314:PHE:CG	1:A:315:ILE:N	2.89	0.40
1:D:64:GLY:HA3	2:D:401:PLR:H5A2	2.03	0.40
1:D:271:LEU:HB2	1:D:333:VAL:HG13	2.03	0.40
1:D:304:LEU:C	1:D:306:ALA:N	2.78	0.40
1:B:128:LYS:HG3	1:C:315:ILE:HD13	2.03	0.40
1:B:278:LEU:HD11	1:B:341:ARG:NH1	2.37	0.40
1:B:314:PHE:C	1:B:316:GLY:N	2.80	0.40
1:C:316:GLY:C	1:C:318:GLY:H	2.28	0.40
1:D:89:HIS:NE2	2:D:401:PLR:H4A2	2.36	0.40
1:D:89:HIS:CE1	2:D:401:PLR:H5A1	2.57	0.40
1:B:81:SER:HB3	1:B:135:PRO:HB2	2.03	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305:ARG:NH2	1:C:346:ALA:O[4_548]	2.11	0.09
1:A:305:ARG:HH22	1:C:346:ALA:O[4_548]	1.59	0.01

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	345/346 (100%)	320 (93%)	20 (6%)	5 (1%)	9	8
1	B	346/346 (100%)	320 (92%)	24 (7%)	2 (1%)	21	25
1	C	342/346 (99%)	322 (94%)	16 (5%)	4 (1%)	10	10
1	D	342/346 (99%)	324 (95%)	14 (4%)	4 (1%)	10	10
All	All	1375/1384 (99%)	1286 (94%)	74 (5%)	15 (1%)	11	11

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	305	ARG
1	C	298	GLU
1	D	298	GLU
1	D	304	LEU
1	D	305	ARG
1	A	318	GLY
1	D	306	ALA
1	C	129	ARG
1	B	317	SER
1	A	304	LEU
1	A	315	ILE
1	C	289	ALA
1	C	317	SER
1	B	315	ILE
1	A	275	VAL

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	276/277 (100%)	275 (100%)	1 (0%)	84	91
1	B	277/277 (100%)	274 (99%)	3 (1%)	65	79
1	C	273/277 (99%)	268 (98%)	5 (2%)	51	68
1	D	273/277 (99%)	266 (97%)	7 (3%)	40	56
All	All	1099/1108 (99%)	1083 (98%)	16 (2%)	57	72

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	344	ILE
1	B	103	ASN
1	B	112	SER
1	B	200	VAL
1	C	113	VAL
1	C	192	LEU
1	C	292	VAL
1	C	313	THR
1	C	344	ILE
1	D	42	GLU
1	D	112	SER
1	D	113	VAL
1	D	125	VAL
1	D	301	ILE
1	D	304	LEU
1	D	321	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	103	ASN
1	A	170	ASN
1	C	77	GLN
1	D	103	ASN
1	D	120	GLN
1	D	180	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

8 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
2	PLR	D	401	-	15,15,15	2.03	7 (46%)	21,22,22	2.02	6 (28%)
2	PLR	C	401	-	15,15,15	2.20	6 (40%)	21,22,22	1.63	4 (19%)
2	PLR	B	401	-	15,15,15	2.21	7 (46%)	21,22,22	2.22	10 (47%)
3	GOL	A	402	-	5,5,5	0.49	0	5,5,5	1.26	1 (20%)
3	GOL	C	402	-	5,5,5	0.64	0	5,5,5	0.90	0
3	GOL	D	402	-	5,5,5	0.50	0	5,5,5	0.65	0
2	PLR	A	401	-	15,15,15	2.24	7 (46%)	21,22,22	2.23	8 (38%)
3	GOL	B	402	-	5,5,5	0.70	0	5,5,5	1.21	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
2	PLR	D	401	-	-	5/6/6/6	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PLR	C	401	-	-	5/6/6/6	0/1/1/1
2	PLR	B	401	-	-	3/6/6/6	0/1/1/1
3	GOL	A	402	-	-	2/4/4/4	-
3	GOL	C	402	-	-	2/4/4/4	-
3	GOL	D	402	-	-	0/4/4/4	-
2	PLR	A	401	-	-	3/6/6/6	0/1/1/1
3	GOL	B	402	-	-	2/4/4/4	-

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	401	PLR	P-O1P	4.36	1.64	1.50
2	A	401	PLR	P-O1P	4.32	1.63	1.50
2	B	401	PLR	P-O4P	3.82	1.72	1.60
2	A	401	PLR	C2-N1	3.34	1.39	1.33
2	B	401	PLR	C2-N1	3.34	1.39	1.33
2	D	401	PLR	C2-N1	3.15	1.39	1.33
2	A	401	PLR	C4A-C4	3.10	1.57	1.51
2	B	401	PLR	O3-C3	3.04	1.43	1.36
2	C	401	PLR	C4A-C4	2.98	1.57	1.51
2	B	401	PLR	C2A-C2	2.97	1.55	1.50
2	D	401	PLR	P-O4P	2.96	1.69	1.60
2	A	401	PLR	C2A-C2	2.95	1.55	1.50
2	C	401	PLR	C2-N1	2.92	1.39	1.33
2	C	401	PLR	P-O4P	2.90	1.69	1.60
2	D	401	PLR	C2A-C2	2.81	1.54	1.50
2	C	401	PLR	O3-C3	2.77	1.43	1.36
2	A	401	PLR	P-O4P	2.75	1.69	1.60
2	D	401	PLR	C4A-C4	2.51	1.56	1.51
2	A	401	PLR	C3-C2	2.48	1.43	1.41
2	D	401	PLR	P-O3P	2.43	1.63	1.54
2	D	401	PLR	O3-C3	2.41	1.42	1.36
2	B	401	PLR	P-O3P	2.40	1.63	1.54
2	A	401	PLR	O3-C3	2.30	1.42	1.36
2	D	401	PLR	P-O1P	2.27	1.57	1.50
2	C	401	PLR	C6-C5	2.20	1.42	1.37
2	B	401	PLR	C4A-C4	2.14	1.55	1.51
2	B	401	PLR	C6-C5	2.09	1.41	1.37

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	PLR	C6-C5-C4	6.16	123.15	118.10
2	D	401	PLR	C6-C5-C4	5.02	122.21	118.10
2	D	401	PLR	O2P-P-O4P	4.51	118.42	106.67
2	C	401	PLR	C6-C5-C4	4.44	121.74	118.10
2	B	401	PLR	O2P-P-O4P	4.37	118.08	106.67
2	B	401	PLR	C2A-C2-C3	-3.83	116.32	120.80
2	D	401	PLR	C5-C6-N1	-3.48	118.17	123.83
2	B	401	PLR	C4A-C4-C5	3.37	124.42	120.94
2	C	401	PLR	C2A-C2-C3	-3.14	117.12	120.80
2	B	401	PLR	C2A-C2-N1	3.08	123.45	117.64
2	B	401	PLR	C6-C5-C4	3.08	120.62	118.10
2	B	401	PLR	O3P-P-O4P	-3.03	98.77	106.67
2	A	401	PLR	O2P-P-O4P	3.02	114.54	106.67
2	A	401	PLR	C4A-C4-C5	2.96	123.99	120.94
2	B	401	PLR	O4P-C5A-C5	2.92	114.82	109.36
2	A	401	PLR	C2A-C2-C3	-2.77	117.55	120.80
2	A	401	PLR	C5-C6-N1	-2.76	119.35	123.83
2	A	401	PLR	C2A-C2-N1	2.60	122.54	117.64
2	A	401	PLR	O2P-P-O1P	2.59	120.94	110.83
2	D	401	PLR	O4P-P-O1P	-2.57	99.50	106.44
2	B	401	PLR	C4A-C4-C3	-2.45	116.44	120.52
2	D	401	PLR	O3P-P-O4P	-2.42	100.35	106.67
2	B	401	PLR	O3P-P-O2P	2.14	115.82	107.80
2	B	401	PLR	C5-C6-N1	-2.10	120.41	123.83
2	A	401	PLR	O3P-P-O4P	-2.10	101.20	106.67
2	C	401	PLR	C5A-C5-C4	-2.09	118.51	122.64
2	D	401	PLR	O3P-P-O2P	2.06	115.51	107.80
2	C	401	PLR	O2P-P-O1P	2.02	118.72	110.83
3	A	402	GOL	O2-C2-C3	2.00	117.47	109.18

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	PLR	C5A-O4P-P-O2P
2	A	401	PLR	C5A-O4P-P-O3P
2	B	401	PLR	C5A-O4P-P-O2P
2	B	401	PLR	C5A-O4P-P-O3P
2	C	401	PLR	C5A-O4P-P-O1P
2	C	401	PLR	C5A-O4P-P-O2P
2	C	401	PLR	C5A-O4P-P-O3P
2	D	401	PLR	C4-C5-C5A-O4P
2	D	401	PLR	C6-C5-C5A-O4P

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Mol	Chain	Res	Type	Atoms
2	D	401	PLR	C5A-O4P-P-O2P
2	D	401	PLR	C5A-O4P-P-O3P
3	A	402	GOL	C1-C2-C3-O3
3	B	402	GOL	C1-C2-C3-O3
3	C	402	GOL	O1-C1-C2-C3
3	A	402	GOL	O2-C2-C3-O3
3	B	402	GOL	O2-C2-C3-O3
3	C	402	GOL	O1-C1-C2-O2
2	B	401	PLR	C5A-O4P-P-O1P
2	D	401	PLR	C5A-O4P-P-O1P
2	C	401	PLR	C6-C5-C5A-O4P
2	C	401	PLR	C4-C5-C5A-O4P
2	A	401	PLR	C5A-O4P-P-O1P

There are no ring outliers.

7 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	401	PLR	6	0
2	C	401	PLR	8	0
2	B	401	PLR	4	0
3	A	402	GOL	1	0
3	C	402	GOL	2	0
2	A	401	PLR	3	0
3	B	402	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 [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	343/346 (99%)	0.18	27 (7%)	18	19	21, 51, 99, 132	11 (3%)
1	B	344/346 (99%)	0.18	25 (7%)	21	22	19, 54, 96, 120	8 (2%)
1	C	342/346 (98%)	0.22	25 (7%)	21	22	22, 54, 93, 133	9 (2%)
1	D	342/346 (98%)	0.23	31 (9%)	15	15	22, 54, 97, 114	11 (3%)
All	All	1371/1384 (99%)	0.21	108 (7%)	18	19	19, 53, 97, 133	39 (2%)

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	306	ALA	6.4
1	A	300	ALA	6.2
1	C	5	SER	5.8
1	D	34	ALA	5.2
1	A	316	GLY	5.2
1	D	31	HIS	5.1
1	A	317	SER	5.1
1	A	314	PHE	4.8
1	A	184	PHE	4.5
1	A	134	TYR	4.5
1	B	127	LEU	4.3
1	A	305	ARG	4.2
1	B	299	HIS	4.2
1	B	34	ALA	4.0
1	D	346	ALA	4.0
1	C	34	ALA	4.0
1	A	315	ILE	3.9
1	A	304	LEU	3.9
1	C	31	HIS	3.9
1	C	129	ARG	3.8
1	B	314	PHE	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	299	HIS	3.7
1	D	30	GLY	3.7
1	A	306	ALA	3.7
1	C	32	ASP	3.7
1	C	184	PHE	3.6
1	B	306	ALA	3.6
1	A	299	HIS	3.6
1	B	317	SER	3.5
1	B	307	LYS	3.5
1	B	346	ALA	3.5
1	D	300	ALA	3.5
1	B	184	PHE	3.4
1	D	314	PHE	3.4
1	B	31	HIS	3.4
1	C	316	GLY	3.4
1	B	302	GLU	3.3
1	D	317	SER	3.3
1	B	134	TYR	3.3
1	D	5	SER	3.3
1	C	134	TYR	3.2
1	C	30	GLY	3.2
1	B	315	ILE	3.2
1	D	305	ARG	3.2
1	D	134	TYR	3.2
1	D	304	LEU	3.2
1	C	127	LEU	3.1
1	D	35	TYR	3.1
1	A	4	LYS	3.1
1	B	295	GLN	3.1
1	D	315	ILE	3.0
1	D	184	PHE	3.0
1	A	303	ALA	3.0
1	C	299	HIS	3.0
1	D	6	GLN	3.0
1	A	346	ALA	3.0
1	A	130	GLN	3.0
1	C	304	LEU	2.9
1	B	300	ALA	2.9
1	A	31	HIS	2.8
1	A	343	ILE	2.8
1	A	128	LYS	2.8
1	C	315	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	318	GLY	2.8
1	C	35	TYR	2.8
1	C	298	GLU	2.8
1	D	36	GLY	2.7
1	A	129	ARG	2.7
1	C	33	ARG	2.7
1	B	303	ALA	2.7
1	D	303	ALA	2.7
1	C	346	ALA	2.6
1	D	129	ARG	2.6
1	C	305	ARG	2.6
1	A	273	LEU	2.5
1	A	6	GLN	2.5
1	D	7	GLN	2.5
1	D	311	PHE	2.5
1	B	35	TYR	2.4
1	C	123	ARG	2.4
1	D	316	GLY	2.4
1	D	32	ASP	2.4
1	A	318	GLY	2.4
1	D	280	GLY	2.4
1	A	308	GLY	2.3
1	C	300	ALA	2.3
1	B	316	GLY	2.3
1	A	13	TYR	2.3
1	C	301	ILE	2.3
1	C	167	LEU	2.2
1	C	303	ALA	2.2
1	B	305	ARG	2.2
1	B	33	ARG	2.2
1	A	133	HIS	2.2
1	B	3	ASP	2.2
1	B	6	GLN	2.2
1	D	33	ARG	2.1
1	D	13	TYR	2.1
1	D	312	TYR	2.1
1	B	301	ILE	2.1
1	C	314	PHE	2.1
1	D	344	ILE	2.1
1	D	130	GLN	2.1
1	B	298	GLU	2.1
1	A	30	GLY	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	32	ASP	2.0
1	B	32	ASP	2.0
1	D	12	ASN	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(Å ²)	Q<0.9
2	PLR	B	401	15/15	0.58	0.24	46,57,72,72	0
2	PLR	D	401	15/15	0.60	0.23	48,61,76,76	0
2	PLR	C	401	15/15	0.64	0.23	45,58,73,73	0
2	PLR	A	401	15/15	0.67	0.22	45,59,74,74	0
3	GOL	D	402	6/6	0.90	0.19	39,61,71,73	0
3	GOL	B	402	6/6	0.93	0.17	44,57,69,70	0
3	GOL	C	402	6/6	0.94	0.12	35,58,77,80	0
3	GOL	A	402	6/6	0.96	0.10	35,51,66,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.