Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 05:20 pm BST

PDB ID	:	6HTE
Title	:	Sulfolobus solfataricus Tryptophan Synthase B2a
Authors	:	Fleming, J.; Mayans, O.; Bucher, R.
Deposited on	:	2018-10-04
Resolution	:	1.96 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 following versions of software and data (see references ⓘ) were used in the production of this report:

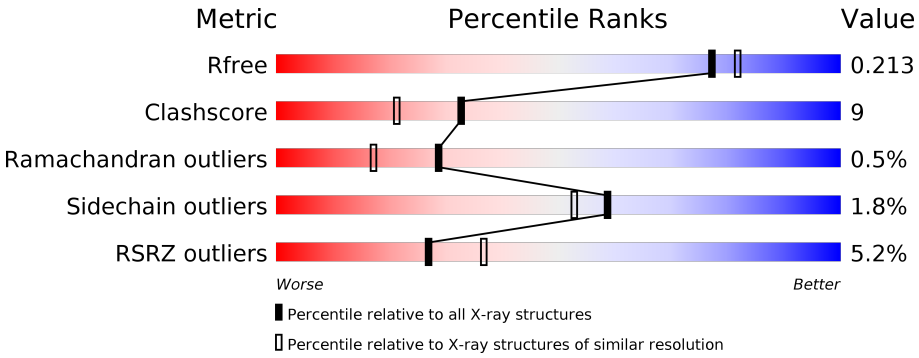
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 1.96 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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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 on 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	B	425	7% 79% 15% • 5%
1	D	425	3% 84% 12% ••

2 Entry composition

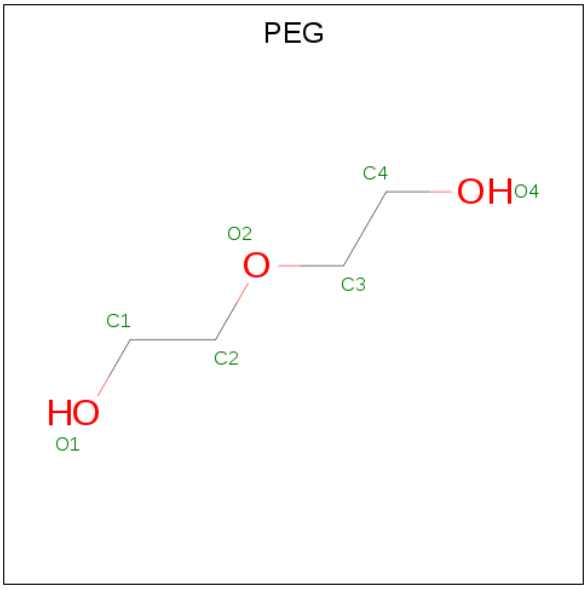
There are 4 unique types of molecules in this entry. The entry contains 13380 atoms, of which 6498 are hydrogens and 0 are deuteriums.

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 Tryptophan synthase beta chain 2.

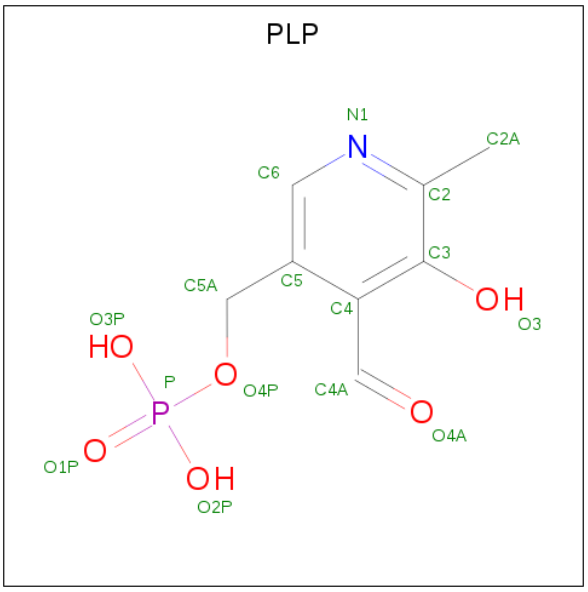
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	405	Total	C	H	N	O	S	0	2	0
			6370	2055	3188	523	592	12			
1	D	415	Total	C	H	N	O	S	0	0	0
			6549	2108	3284	540	605	12			

- Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	H	O	0	0
			17	4	10	3		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total	C	H	N	O	P	0	0
			23	8	8	1	5	1		
3	D	1	Total	C	H	N	O	P	0	0
			23	8	8	1	5	1		

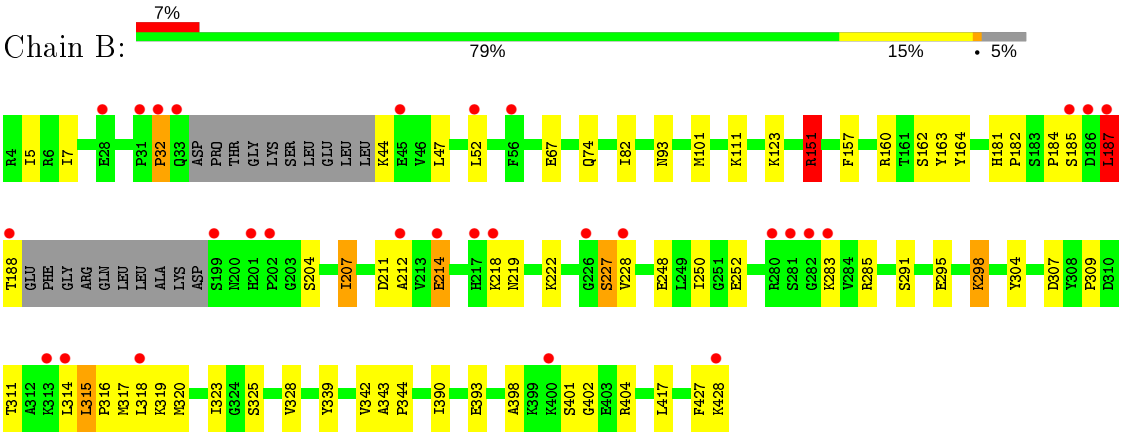
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	191	Total	O	0	0
			191	191		
4	D	207	Total	O	0	0
			207	207		

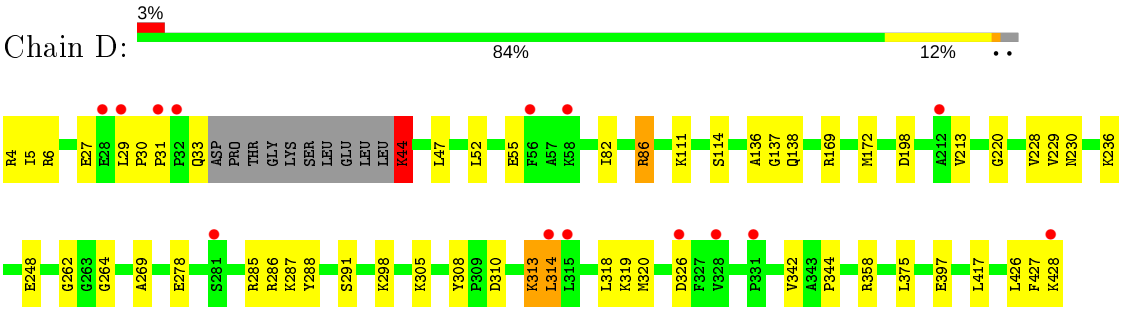
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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: Tryptophan synthase beta chain 2



- Molecule 1: Tryptophan synthase beta chain 2



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	55.89Å 61.71Å 110.36Å 90.00° 99.33° 90.00°	Depositor
Resolution (Å)	19.95 – 1.96 19.95 – 1.96	Depositor EDS
% Data completeness (in resolution range)	99.5 (19.95-1.96) 99.5 (19.95-1.96)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.77 (at 1.96Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.163 , 0.213 0.163 , 0.213	Depositor DCC
R_{free} test set	2655 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	24.6	Xtriage
Anisotropy	0.242	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 51.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	13380	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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: PEG, PLP

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	B	0.53	0/3268	0.64	2/4425 (0.0%)
1	D	0.59	0/3346	0.65	1/4527 (0.0%)
All	All	0.56	0/6614	0.64	3/8952 (0.0%)

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	2
1	D	0	1
All	All	0	3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	151	ARG	NE-CZ-NH2	-7.79	116.40	120.30
1	D	86	ARG	NE-CZ-NH2	-5.84	117.38	120.30
1	B	151	ARG	NE-CZ-NH1	5.78	123.19	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	187	LEU	Peptide
1	B	227	SER	Peptide
1	D	44	LYS	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	B	3182	3188	3178	66	0
1	D	3265	3284	3286	48	0
2	B	7	10	10	3	0
3	B	15	8	7	1	0
3	D	15	8	7	1	0
4	B	191	0	0	14	0
4	D	207	0	0	11	1
All	All	6882	6498	6488	113	1

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 9.

All (113) 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:B:93:ASN:O	1:B:404:ARG:NH1	2.08	0.87
1:D:47:LEU:HD12	1:D:52:LEU:HD21	1.59	0.85
1:B:295:GLU:OE2	4:B:601:HOH:O	1.95	0.83
1:B:5:ILE:HG22	1:B:82:ILE:HG23	1.63	0.80
1:B:44:LYS:N	4:B:607:HOH:O	2.14	0.80
1:B:393:GLU:OE2	4:B:602:HOH:O	1.99	0.80
3:B:502:PLP:O2P	4:B:604:HOH:O	2.02	0.77
1:B:74[B]:GLN:OE1	4:B:603:HOH:O	2.01	0.76
1:B:252:GLU:OE2	4:B:605:HOH:O	2.04	0.75
1:D:33:GLN:NE2	4:D:603:HOH:O	2.08	0.75
1:D:198:ASP:OD2	4:D:601:HOH:O	2.06	0.73
1:D:428:LYS:O	4:D:602:HOH:O	2.07	0.72
1:B:74[B]:GLN:NE2	4:B:609:HOH:O	2.20	0.71
1:B:47:LEU:HD23	1:B:52:LEU:CD2	2.21	0.71
1:B:311:THR:HG22	1:B:325:SER:O	1.92	0.69
1:B:325:SER:OG	4:B:606:HOH:O	2.11	0.69
1:D:426:LEU:O	4:D:604:HOH:O	2.11	0.68
1:B:47:LEU:HD23	1:B:52:LEU:HD21	1.77	0.65
1:D:305:LYS:NZ	4:D:607:HOH:O	2.29	0.65
1:D:47:LEU:CD1	1:D:52:LEU:HD21	2.28	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:MET:C	1:B:318:LEU:HD12	2.19	0.63
1:D:313:LYS:O	1:D:314:LEU:HB3	1.99	0.63
1:B:123:LYS:HD3	2:B:501:PEG:H22	1.80	0.62
1:B:160:ARG:NH2	1:B:184:PRO:O	2.25	0.61
1:D:310:ASP:OD1	1:D:313:LYS:O	2.17	0.61
1:D:287:LYS:NZ	1:D:397:GLU:OE2	2.27	0.61
1:D:262:GLY:O	1:D:298:LYS:HE3	2.01	0.60
1:B:185:SER:O	1:B:188:THR:HG21	2.02	0.60
1:D:55:GLU:OE2	1:D:230:ASN:N	2.36	0.59
1:B:320:MET:HB2	1:B:342:VAL:O	2.01	0.59
1:B:123:LYS:HD3	2:B:501:PEG:C2	2.33	0.58
1:D:114:SER:HB2	1:D:236:LYS:HE3	1.87	0.57
1:B:7:ILE:HD11	1:B:250:ILE:HD13	1.87	0.57
1:B:314:LEU:O	1:B:315:LEU:HB2	2.05	0.57
1:D:213:VAL:HG21	1:D:314:LEU:HD22	1.86	0.57
1:B:182:PRO:O	1:B:185:SER:HB3	2.04	0.56
1:B:328:VAL:HG13	4:B:716:HOH:O	2.06	0.56
1:B:218:LYS:HG3	1:B:219:ASN:OD1	2.06	0.55
1:D:285:ARG:HG3	1:D:286:ARG:N	2.21	0.54
1:B:309:PRO:CG	1:B:315:LEU:HD23	2.37	0.54
1:B:401:SER:N	1:B:402:GLY:HA2	2.23	0.54
1:B:398:ALA:O	1:B:402:GLY:HA2	2.07	0.54
1:D:278:GLU:OE2	1:D:288:TYR:OH	2.20	0.53
1:D:427:PHE:O	1:D:428:LYS:HB3	2.09	0.53
1:B:298:LYS:HA	1:B:323:ILE:HG23	1.90	0.53
1:B:228:VAL:CG2	1:B:343:ALA:HB2	2.38	0.52
1:B:123:LYS:HE2	2:B:501:PEG:H42	1.91	0.52
1:B:307:ASP:OD1	4:B:606:HOH:O	2.19	0.51
1:B:304:TYR:CE1	1:B:319:LYS:HG3	2.45	0.51
1:D:285:ARG:NH2	4:D:614:HOH:O	2.43	0.51
1:B:317:MET:O	1:B:318:LEU:HD12	2.11	0.51
1:B:187:LEU:N	1:B:187:LEU:HD23	2.26	0.51
1:D:55:GLU:OE2	1:D:229:VAL:HB	2.11	0.51
1:B:74[A]:GLN:NE2	1:D:86:ARG:HD3	2.25	0.50
1:D:318:LEU:HB3	1:D:344:PRO:HG3	1.92	0.50
1:D:313:LYS:N	1:D:313:LYS:HD3	2.25	0.50
1:D:30:PRO:CG	1:D:319:LYS:HB3	2.40	0.50
1:D:30:PRO:HG3	1:D:319:LYS:HB3	1.94	0.50
1:D:44:LYS:CD	1:D:44:LYS:N	2.74	0.50
1:B:204:SER:O	1:B:207:ILE:HG23	2.12	0.50
1:B:82:ILE:HD12	1:B:101:MET:HG3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:86:ARG:NE	1:D:375:LEU:O	2.45	0.49
1:B:7:ILE:CD1	1:B:250:ILE:HD13	2.42	0.48
1:D:172:MET:HE1	1:D:417:LEU:HD13	1.96	0.48
1:B:157:PHE:CD2	1:B:212:ALA:HB2	2.48	0.47
1:D:248:GLU:HG3	4:D:694:HOH:O	2.13	0.47
1:B:184:PRO:HB3	1:B:207:ILE:HD11	1.95	0.47
1:D:44:LYS:HD3	1:D:44:LYS:N	2.29	0.47
1:D:111:LYS:NZ	3:D:501:PLP:O3	2.47	0.47
1:B:309:PRO:O	1:B:339:TYR:OH	2.17	0.47
1:B:417:LEU:HD23	4:B:712:HOH:O	2.15	0.47
1:B:248:GLU:HG3	4:B:695:HOH:O	2.15	0.46
1:D:313:LYS:O	1:D:314:LEU:CB	2.63	0.46
1:D:172:MET:HE1	1:D:417:LEU:CD1	2.46	0.46
1:B:291:SER:HB3	1:B:390:ILE:HD11	1.98	0.45
1:D:291:SER:HA	1:D:358:ARG:O	2.17	0.45
1:B:309:PRO:HG2	1:B:315:LEU:HD23	1.98	0.45
1:B:162:SER:OG	4:B:608:HOH:O	2.20	0.45
1:B:283:LYS:HE3	1:B:285:ARG:HB2	1.98	0.44
1:B:67:GLU:HG2	4:B:729:HOH:O	2.17	0.44
1:B:163:TYR:CE2	1:B:182:PRO:HD3	2.53	0.44
1:B:204:SER:H	1:B:207:ILE:CG2	2.31	0.44
1:D:308:TYR:CE2	1:D:313:LYS:HE2	2.54	0.43
1:B:318:LEU:HD23	1:B:344:PRO:HG2	2.01	0.43
1:D:264:GLY:HA2	4:D:620:HOH:O	2.18	0.43
1:D:320:MET:HB2	1:D:342:VAL:O	2.17	0.43
1:B:222:LYS:HD3	1:B:222:LYS:HA	1.85	0.43
1:D:308:TYR:CE1	1:D:313:LYS:HG3	2.54	0.43
1:B:181:HIS:CE1	1:B:187:LEU:HD11	2.54	0.43
1:D:427:PHE:O	1:D:428:LYS:CB	2.66	0.43
1:B:188:THR:HA	1:B:214:GLU:OE2	2.19	0.43
1:D:220:GLY:N	4:D:619:HOH:O	2.45	0.43
1:D:236:LYS:O	1:D:269:ALA:HB1	2.19	0.43
1:B:228:VAL:HG22	1:B:343:ALA:CB	2.50	0.42
1:D:220:GLY:HA2	4:D:619:HOH:O	2.19	0.42
1:B:151:ARG:HD3	1:D:86:ARG:HH22	1.83	0.42
1:D:136:ALA:HB3	1:D:138:GLN:OE1	2.19	0.42
1:B:314:LEU:O	1:B:315:LEU:CB	2.68	0.42
1:B:160:ARG:HG3	1:B:164:TYR:HE2	1.84	0.42
1:B:427:PHE:O	1:B:428:LYS:HB2	2.20	0.42
1:B:228:VAL:HG22	1:B:228:VAL:O	2.20	0.41
1:B:207:ILE:O	1:B:211:ASP:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:326:ASP:OD1	1:D:326:ASP:N	2.53	0.41
1:B:184:PRO:HA	1:B:211:ASP:OD2	2.21	0.41
1:B:32:PRO:HB3	1:B:316:PRO:HB2	2.02	0.41
1:B:47:LEU:CD2	1:B:52:LEU:HD21	2.48	0.41
1:D:30:PRO:HA	1:D:31:PRO:HD2	1.95	0.41
1:B:318:LEU:CD2	1:B:344:PRO:HG2	2.51	0.41
1:D:137:GLY:C	1:D:169:ARG:HD3	2.41	0.41
1:D:6:ARG:HD3	4:D:708:HOH:O	2.20	0.41
1:B:228:VAL:HG22	1:B:343:ALA:HB2	2.02	0.40
1:D:5:ILE:HD12	1:D:82:ILE:HG23	2.01	0.40
1:B:160:ARG:HG3	1:B:164:TYR:CE2	2.55	0.40

All (1) 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)
4:D:806:HOH:O	4:D:807:HOH:O[2_555]	2.18	0.02

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	B	401/425 (94%)	383 (96%)	16 (4%)	2 (0%)	29	17
1	D	411/425 (97%)	398 (97%)	11 (3%)	2 (0%)	29	17
All	All	812/850 (96%)	781 (96%)	27 (3%)	4 (0%)	29	17

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	314	LEU
1	B	315	LEU

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Mol	Chain	Res	Type
1	D	228	VAL
1	B	32	PRO

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	B	339/356 (95%)	332 (98%)	7 (2%)	53	46
1	D	347/356 (98%)	342 (99%)	5 (1%)	67	62
All	All	686/712 (96%)	674 (98%)	12 (2%)	59	55

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	111	LYS
1	B	151	ARG
1	B	187	LEU
1	B	207	ILE
1	B	214	GLU
1	B	227	SER
1	B	298	LYS
1	D	4	ARG
1	D	27	GLU
1	D	29	LEU
1	D	44	LYS
1	D	313	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

3 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	PEG	B	501	-	6,6,6	0.50	0	5,5,5	1.07	0
3	PLP	D	501	1	15,15,16	2.37	7 (46%)	20,22,23	1.74	3 (15%)
3	PLP	B	502	1	15,15,16	2.05	4 (26%)	20,22,23	1.66	4 (20%)

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	PEG	B	501	-	-	4/4/4/4	-
3	PLP	D	501	1	-	2/6/6/8	0/1/1/1
3	PLP	B	502	1	-	3/6/6/8	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	501	PLP	C4A-C4	5.07	1.62	1.51
3	B	502	PLP	C4A-C4	4.77	1.61	1.51
3	D	501	PLP	C5-C4	-3.71	1.36	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	501	PLP	C2A-C2	3.33	1.56	1.50
3	B	502	PLP	C5-C4	-3.10	1.37	1.40
3	B	502	PLP	C2A-C2	2.81	1.55	1.50
3	D	501	PLP	C6-N1	2.76	1.40	1.34
3	B	502	PLP	C5A-C5	2.46	1.57	1.50
3	D	501	PLP	P-O4P	-2.33	1.52	1.60
3	D	501	PLP	P-O3P	-2.15	1.46	1.54
3	D	501	PLP	C3-C2	2.07	1.43	1.40

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	501	PLP	O4P-C5A-C5	-4.61	100.56	109.35
3	B	502	PLP	C4A-C4-C5	4.31	125.37	120.94
3	D	501	PLP	C6-C5-C4	3.78	121.13	118.16
3	B	502	PLP	C4A-C4-C3	-2.93	115.53	120.50
3	D	501	PLP	C5-C6-N1	-2.91	118.96	123.82
3	B	502	PLP	O3P-P-O4P	2.62	113.70	106.73
3	B	502	PLP	C5A-C5-C6	-2.30	115.59	119.37

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	501	PLP	C4-C5-C5A-O4P
3	B	502	PLP	C5A-O4P-P-O1P
3	B	502	PLP	C5A-O4P-P-O2P
3	B	502	PLP	C5A-O4P-P-O3P
2	B	501	PEG	O2-C3-C4-O4
2	B	501	PEG	C4-C3-O2-C2
3	D	501	PLP	C6-C5-C5A-O4P
2	B	501	PEG	C1-C2-O2-C3
2	B	501	PEG	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	PEG	3	0
3	D	501	PLP	1	0
3	B	502	PLP	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	B	405/425 (95%)	0.23	29 (7%)	15	23	12, 28, 60, 81	0
1	D	415/425 (97%)	-0.09	14 (3%)	45	55	11, 25, 51, 79	0
All	All	820/850 (96%)	0.07	43 (5%)	27	37	11, 26, 55, 81	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	188	THR	8.9
1	B	199	SER	6.6
1	B	187	LEU	5.7
1	B	56	PHE	5.1
1	B	186	ASP	4.9
1	D	28	GLU	4.7
1	B	185	SER	4.0
1	B	33	GLN	3.8
1	D	56	PHE	3.7
1	D	428	LYS	3.7
1	B	226	GLY	3.4
1	B	202	PRO	3.2
1	B	31	PRO	3.2
1	D	31	PRO	3.1
1	D	32	PRO	3.0
1	B	282	GLY	3.0
1	B	217	HIS	2.9
1	B	314	LEU	2.9
1	D	314	LEU	2.9
1	B	313	LYS	2.8
1	B	214	GLU	2.8
1	B	52	LEU	2.7
1	B	318	LEU	2.7
1	B	283	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	328	VAL	2.7
1	D	326	ASP	2.7
1	B	201	HIS	2.5
1	D	281	SER	2.5
1	B	218	LYS	2.5
1	D	29	LEU	2.5
1	B	45	GLU	2.4
1	D	315	LEU	2.4
1	B	400	LYS	2.3
1	B	32	PRO	2.3
1	B	28	GLU	2.3
1	B	212	ALA	2.2
1	B	281	SER	2.2
1	D	331	PRO	2.1
1	B	428	LYS	2.1
1	B	228	VAL	2.1
1	B	280	ARG	2.1
1	D	212	ALA	2.1
1	D	58	LYS	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 carbohydrates 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	PEG	B	501	7/7	0.90	0.40	39,54,66,70	0
3	PLP	D	501	15/16	0.98	0.08	15,22,31,35	0
3	PLP	B	502	15/16	0.98	0.07	17,23,32,37	0

6.5 Other polymers (i)

There are no such residues in this entry.