



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:21 PM UTC

PDB ID : 4LW4 / pdb_00004lw4
Title : Structural changes during cysteine desulfurase CsdA and sulfur-acceptor CsdE interactions provide insight into the trans-persulfuration
Authors : Kim, S.; Park, S.Y.
Deposited on : 2013-07-26
Resolution : 2.01 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

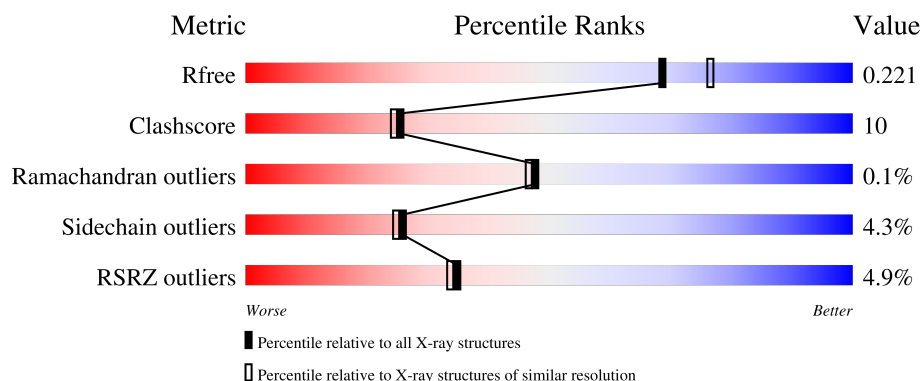
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	404	
1	B	404	
2	C	150	
2	D	150	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8926 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 Cysteine sulfinate desulfinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	392	Total	C	N	O	S	0	1	0
			2976	1892	512	560	12			
1	B	398	Total	C	N	O	S	0	7	0
			3065	1947	530	576	12			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q46925
A	-1	SER	-	expression tag	UNP Q46925
A	0	HIS	-	expression tag	UNP Q46925
B	-2	GLY	-	expression tag	UNP Q46925
B	-1	SER	-	expression tag	UNP Q46925
B	0	HIS	-	expression tag	UNP Q46925

- Molecule 2 is a protein called Cysteine desulfuration protein CsdE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	145	Total	C	N	O	S	0	3	0
			1123	712	198	210	3			
2	D	137	Total	C	N	O	S	0	1	0
			1047	665	183	196	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	GLY	-	expression tag	UNP D5CZ88
C	-1	SER	-	expression tag	UNP D5CZ88
C	0	HIS	-	expression tag	UNP D5CZ88
C	49	ASP	GLU	engineered mutation	UNP D5CZ88
D	-2	GLY	-	expression tag	UNP D5CZ88
D	-1	SER	-	expression tag	UNP D5CZ88

Continued on next page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	0	HIS	-	expression tag	UNP D5CZ88
D	49	ASP	GLU	engineered mutation	UNP D5CZ88

- PLP
-
- The diagram shows the chemical structure of Pyridoxal Phosphate (PLP). It consists of a pyridine ring with a phosphate group at the 3-position and an aldehyde group at the 4-position. The pyridine ring is colored blue, with the nitrogen atom labeled N1. The carbon atoms are labeled C2, C3, C4, C5, and C6. The phosphate group is colored purple, with the phosphorus atom labeled P. The oxygen atoms are labeled O1P, O2P, O3P, O4P, and O5P. The aldehyde group is colored red, with the carbonyl oxygen labeled O4A and the hydroxyl group labeled O3. The hydroxyl group at the 3-position is colored red, with the oxygen labeled O3 and the hydrogen labeled H3. The hydroxyl group at the 4-position is colored red, with the oxygen labeled O4A and the hydrogen labeled H4A. The hydroxyl group at the 5-position is colored red, with the oxygen labeled O5P and the hydrogen labeled H5P. The hydroxyl group at the 6-position is colored red, with the oxygen labeled O6P and the hydrogen labeled H6P.

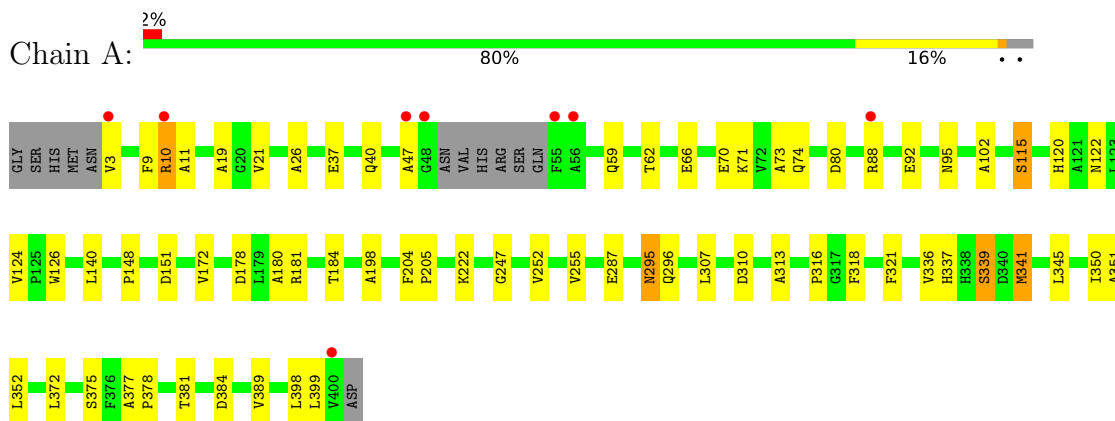
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 15	C 8	N 1	O 5	P 1	0	0
3	B	1	Total 15	C 8	N 1	O 5	P 1	0	0

- | Mol | Chain | Residues | Atoms | ZeroOcc | AltConf |
|-----|-------|----------|--------------------|---------|---------|
| 4 | A | 289 | Total O
289 289 | 0 | 0 |
| 4 | B | 316 | Total O
316 316 | 0 | 0 |
| 4 | C | 74 | Total O
74 74 | 0 | 0 |
| 4 | D | 6 | Total O
6 6 | 0 | 0 |

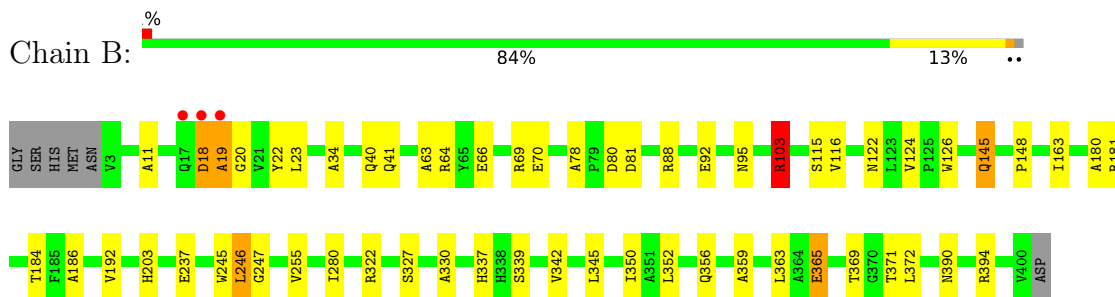
3 Residue-property plots

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.

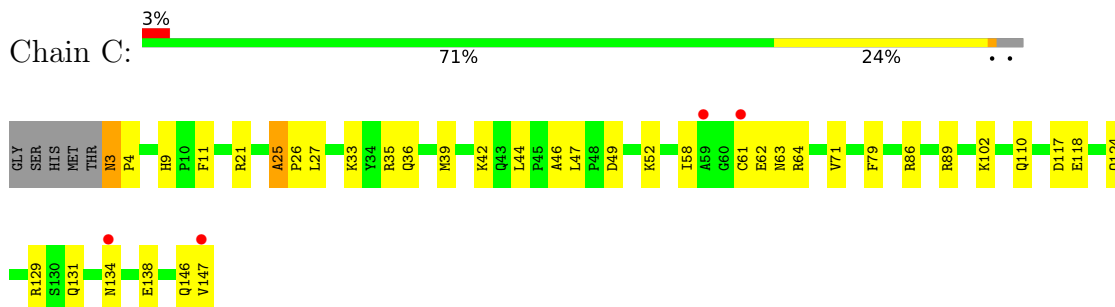
- Molecule 1: Cysteine sulfinate desulfinase



- Molecule 1: Cysteine sulfinate desulfinase

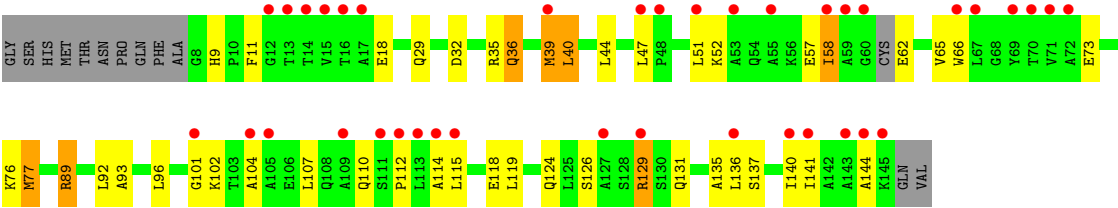


- Molecule 2: Cysteine desulfuration protein CsdE



- Molecule 2: Cysteine desulfuration protein CsdE





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	77.14Å 76.87Å 89.36Å 90.00° 104.48° 90.00°	Depositor
Resolution (Å)	50.00 – 2.01 50.00 – 2.01	Depositor EDS
% Data completeness (in resolution range)	99.4 (50.00-2.01) 99.6 (50.00-2.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.08 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.159 , 0.218 0.163 , 0.221	Depositor DCC
R_{free} test set	3411 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	19.8	Xtriage
Anisotropy	0.256	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8926	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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: 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	A	1.39	13/3044 (0.4%)	1.13	4/4147 (0.1%)
1	B	1.37	8/3153 (0.3%)	1.12	4/4293 (0.1%)
2	C	1.19	1/1151 (0.1%)	1.10	6/1558 (0.4%)
2	D	0.86	0/1066	1.06	2/1440 (0.1%)
All	All	1.30	22/8414 (0.3%)	1.11	16/11438 (0.1%)

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	124	VAL	CA-CB	9.14	1.58	1.54
1	A	351	ALA	CA-CB	-7.39	1.43	1.53
1	A	377	ALA	CA-CB	6.71	1.64	1.53
1	B	186	ALA	CA-CB	6.37	1.63	1.53
1	A	389	VAL	CA-CB	6.34	1.62	1.54
1	A	172	VAL	CA-CB	6.20	1.62	1.54
1	A	21	VAL	C-O	-6.08	1.17	1.24
1	A	350	ILE	C-O	-6.03	1.17	1.24
1	A	19	ALA	C-O	-5.98	1.16	1.24
1	A	321	PHE	C-O	-5.90	1.16	1.24
1	A	10	ARG	C-O	-5.89	1.17	1.24
1	A	124	VAL	CA-C	5.89	1.58	1.52
1	B	322	ARG	C-O	-5.74	1.17	1.23
1	B	11	ALA	C-O	-5.68	1.16	1.24
1	A	310	ASP	N-CA	5.42	1.52	1.46
1	B	359	ALA	CA-CB	5.33	1.60	1.53
1	A	11	ALA	C-O	-5.26	1.17	1.24
2	C	129	ARG	C-O	-5.17	1.16	1.24
1	B	63	ALA	CA-CB	5.14	1.61	1.53
1	A	73	ALA	CA-CB	5.03	1.61	1.53
1	B	103	ARG	CA-C	5.03	1.58	1.52
1	B	34	ALA	CA-CB	-5.00	1.45	1.53

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	78	ALA	CA-C-N	-6.25	113.32	119.64
1	B	78	ALA	C-N-CA	-6.25	113.32	119.64
2	D	73	GLU	CA-C-N	6.18	129.18	120.28
2	D	73	GLU	C-N-CA	6.18	129.18	120.28
1	B	280	ILE	N-CA-C	5.70	115.89	110.42
1	A	399	LEU	N-CA-C	5.68	120.28	113.23
2	C	62	GLU	N-CA-C	5.59	118.31	111.82
2	C	25	ALA	CA-C-N	5.54	125.64	119.87
2	C	25	ALA	C-N-CA	5.54	125.64	119.87
2	C	3	ASN	CA-C-N	-5.41	114.21	119.78
2	C	3	ASN	C-N-CA	-5.41	114.21	119.78
1	A	102	ALA	N-CA-C	5.38	116.95	111.14
1	A	255	VAL	CB-CA-C	-5.29	101.55	110.71
1	A	384	ASP	N-CA-C	-5.11	105.79	111.36
1	B	327	SER	N-CA-C	-5.11	107.00	113.18
2	C	117	ASP	N-CA-C	-5.06	105.84	111.36

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	2976	0	2940	58	0
1	B	3065	0	3041	56	0
2	C	1123	0	1141	26	0
2	D	1047	0	1067	43	0
3	A	15	0	7	1	0
3	B	15	0	6	0	0
4	A	289	0	0	3	0
4	B	316	0	0	11	0
4	C	74	0	0	3	0
4	D	6	0	0	0	0
All	All	8926	0	8202	158	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (158) 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:10:ARG:NH1	1:A:381:THR:HG21	1.28	1.45
1:A:10:ARG:NH1	1:A:381:THR:CG2	1.91	1.33
1:B:66:GLU:HG2	4:B:790:HOH:O	1.42	1.16
2:D:126:SER:O	2:D:129:ARG:HG2	1.62	1.00
1:B:70:GLU:HG3	4:B:908:HOH:O	1.67	0.94
1:A:88:ARG:HH21	1:B:88[A]:ARG:CD	1.82	0.93
1:A:10:ARG:HH12	1:A:381:THR:HG21	1.12	0.91
2:D:77:MET:HE3	2:D:107:LEU:HD12	1.51	0.91
1:A:10:ARG:HH11	1:A:381:THR:HG21	1.17	0.90
1:A:151:ASP:OD1	1:A:181:ARG:NH1	2.04	0.90
1:A:10:ARG:NH1	1:A:381:THR:HG23	1.86	0.88
1:A:10:ARG:HH11	1:A:381:THR:CG2	1.74	0.87
1:B:342:VAL:HG11	2:D:62:GLU:OE2	1.77	0.85
1:A:95:ASN:HD21	1:B:247:GLY:H	1.22	0.84
1:A:88:ARG:HH21	1:B:88[A]:ARG:HD2	1.42	0.84
1:A:337:HIS:HD2	1:A:339:SER:HB2	1.40	0.84
1:A:337:HIS:CD2	1:A:339:SER:HB2	2.13	0.83
1:B:88[B]:ARG:CZ	4:B:792:HOH:O	2.26	0.83
2:C:58:ILE:O	2:C:58:ILE:HG22	1.77	0.82
2:D:126:SER:O	2:D:129:ARG:CG	2.29	0.81
1:A:247:GLY:H	1:B:95:ASN:HD21	1.28	0.80
1:B:369:THR:HG22	4:B:839:HOH:O	1.81	0.79
2:D:40:LEU:HD22	2:D:93:ALA:CB	2.14	0.78
2:C:9:HIS:HD2	2:C:11:PHE:H	1.31	0.77
2:D:40:LEU:HD21	2:D:93:ALA:HB1	1.65	0.77
2:C:21:ARG:HD3	2:C:118:GLU:OE2	1.87	0.74
2:D:44:LEU:HD23	2:D:89:ARG:HD2	1.68	0.74
2:D:40:LEU:CD2	2:D:93:ALA:CB	2.66	0.73
1:B:88[B]:ARG:NH1	4:B:792:HOH:O	2.21	0.73
1:A:62:THR:O	1:A:66[A]:GLU:HG2	1.88	0.72
1:A:88:ARG:HH21	1:B:88[A]:ARG:HD3	1.55	0.71
1:B:356:GLN:OE1	1:B:363:LEU:HD12	1.90	0.71
1:A:180:ALA:O	1:A:184:THR:HG23	1.92	0.70
1:A:316:PRO:HD2	4:A:738:HOH:O	1.92	0.69
1:B:371:THR:HG22	1:B:372:LEU:O	1.92	0.69
2:D:35:ARG:O	2:D:39[A]:MET:HG2	1.92	0.69
2:D:40:LEU:CD2	2:D:93:ALA:HB1	2.23	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:44:LEU:O	2:C:89[B]:ARG:NH1	2.17	0.67
1:A:37:GLU:HB3	4:A:775:HOH:O	1.96	0.65
1:A:339:SER:OG	2:C:63:ASN:HB2	1.95	0.65
1:B:163:ILE:HD13	1:B:192:VAL:HG22	1.79	0.65
2:D:40:LEU:HD22	2:D:93:ALA:HB3	1.79	0.64
1:A:10:ARG:HH12	1:A:381:THR:CG2	1.82	0.64
1:B:40[B]:GLN:HE21	1:B:41:GLN:HE21	1.45	0.63
1:B:337:HIS:CD2	1:B:339:SER:H	2.15	0.63
2:D:40:LEU:C	2:D:40:LEU:HD23	2.23	0.63
1:B:339:SER:HB2	2:D:131:GLN:NE2	2.13	0.63
2:C:47:LEU:O	2:C:52:LYS:HE2	1.98	0.63
2:D:36:GLN:HA	2:D:39[A]:MET:HG3	1.82	0.62
2:C:42:LYS:HZ3	2:C:86:ARG:HH21	1.48	0.62
1:A:95:ASN:ND2	1:B:247:GLY:H	1.95	0.61
1:A:47:ALA:O	1:B:20:GLY:HA3	2.01	0.61
2:C:46:ALA:HA	2:C:89[B]:ARG:NH2	2.16	0.61
2:C:42:LYS:HZ3	2:C:86:ARG:NH2	1.98	0.60
2:D:40:LEU:HD23	2:D:40:LEU:O	2.02	0.60
1:A:337:HIS:CD2	1:A:339:SER:H	2.19	0.59
2:D:47:LEU:HD23	2:D:52:LYS:HG2	1.84	0.59
1:B:181:ARG:HG3	4:B:815:HOH:O	2.01	0.59
2:C:42:LYS:NZ	2:C:86:ARG:NH2	2.51	0.59
2:C:58:ILE:O	2:C:58:ILE:CG2	2.47	0.59
1:B:339:SER:HB2	2:D:131:GLN:HE21	1.68	0.58
1:A:122:ASN:ND2	1:A:126:TRP:HE1	2.00	0.58
1:A:88:ARG:NH2	1:B:88[A]:ARG:CD	2.62	0.58
1:A:47:ALA:HB1	1:B:22:TYR:HB2	1.86	0.57
1:A:88:ARG:NH2	1:B:88[A]:ARG:HD2	2.17	0.57
1:B:337:HIS:HD2	1:B:339:SER:H	1.52	0.57
2:C:9:HIS:HD2	2:C:11:PHE:N	2.03	0.57
2:D:76:LYS:HG2	2:D:101:GLY:O	2.05	0.57
2:C:9:HIS:HE1	2:C:79:PHE:O	1.87	0.56
2:D:57:GLU:HG3	2:D:66:TRP:NE1	2.20	0.56
2:D:9:HIS:HD2	2:D:11:PHE:H	1.52	0.56
1:A:88:ARG:NH1	1:B:246:LEU:CD2	2.69	0.55
2:D:40:LEU:CD2	2:D:40:LEU:C	2.81	0.54
1:A:337:HIS:HD2	1:A:339:SER:H	1.55	0.53
2:D:114:ALA:O	2:D:118:GLU:HG3	2.08	0.53
1:B:245:TRP:CD1	1:B:246:LEU:HD13	2.43	0.53
1:A:88:ARG:NH1	1:B:246:LEU:HD21	2.23	0.53
1:A:247:GLY:H	1:B:95:ASN:ND2	2.01	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:ILE:HD13	1:B:192:VAL:CG2	2.39	0.53
2:D:77:MET:HE2	2:D:104:ALA:N	2.25	0.52
1:B:64:ARG:HG2	4:B:672:HOH:O	2.08	0.52
2:D:77:MET:CE	2:D:107:LEU:HD12	2.34	0.52
1:B:23:LEU:HB2	1:B:352:LEU:HD12	1.91	0.52
2:D:110:GLN:OE1	2:D:112:PRO:HA	2.10	0.52
1:A:295:ASN:HD22	1:A:295:ASN:H	1.58	0.51
1:B:116:VAL:HG21	1:B:365:GLU:HG3	1.92	0.51
1:A:295:ASN:H	1:A:295:ASN:ND2	2.09	0.51
2:D:129:ARG:HB2	2:D:129:ARG:NH1	2.27	0.50
1:A:88:ARG:NH1	1:B:92:GLU:OE2	2.45	0.50
2:D:77:MET:HE2	2:D:104:ALA:CA	2.42	0.50
1:B:18:ASP:O	1:B:19:ALA:O	2.30	0.49
2:C:9:HIS:CE1	2:C:79:PHE:O	2.66	0.49
2:C:131:GLN:O	2:C:134[B]:ASN:OD1	2.30	0.49
1:B:350:ILE:HG22	1:B:352:LEU:HD13	1.95	0.49
2:C:3:ASN:N	2:C:4:PRO:HD2	2.27	0.49
2:C:35:ARG:NH2	4:C:236:HOH:O	2.46	0.49
1:B:122:ASN:ND2	1:B:126:TRP:HE1	2.10	0.49
2:D:58:ILE:HD12	2:D:135:ALA:HB1	1.95	0.49
2:C:27:LEU:HD11	2:C:36:GLN:HG3	1.94	0.48
2:D:77:MET:HG3	2:D:102:LYS:O	2.12	0.48
2:D:92:LEU:HD13	2:D:136:LEU:HD13	1.95	0.48
1:A:88:ARG:CZ	1:B:92:GLU:OE2	2.61	0.47
2:D:115:LEU:O	2:D:119:LEU:HG	2.14	0.47
2:D:29:GLN:HB2	2:D:32:ASP:OD2	2.14	0.47
1:A:88:ARG:HE	1:B:88[A]:ARG:CG	2.28	0.47
1:A:115:SER:HA	1:A:140:LEU:HB3	1.96	0.47
2:D:129:ARG:HB2	2:D:129:ARG:HH11	1.79	0.47
2:D:77:MET:CE	2:D:104:ALA:HA	2.45	0.47
2:C:25:ALA:N	2:C:26:PRO:CD	2.78	0.47
1:A:70:GLU:O	1:A:74:GLN:HG3	2.14	0.47
2:D:52:LYS:HD2	2:D:66:TRP:CE2	2.49	0.47
1:A:88:ARG:NH2	1:A:92:GLU:OE2	2.48	0.46
1:A:352:LEU:C	1:A:352:LEU:HD12	2.40	0.46
1:A:178:ASP:OD1	1:A:178:ASP:C	2.59	0.46
1:B:390:ASN:HD21	1:B:394:ARG:CZ	2.29	0.46
1:A:120:HIS:CD2	1:B:255:VAL:HG11	2.51	0.45
2:D:126:SER:O	2:D:129:ARG:HG3	2.16	0.45
1:A:3:VAL:HA	1:A:296:GLN:OE1	2.17	0.44
1:B:18:ASP:O	1:B:19:ALA:C	2.61	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:PHE:O	1:A:205:PRO:C	2.59	0.44
1:B:180:ALA:O	1:B:184:THR:HG23	2.18	0.44
1:B:64:ARG:NH2	4:B:696:HOH:O	2.51	0.44
1:B:345:LEU:HD13	1:B:352:LEU:HD22	1.99	0.44
1:A:88:ARG:NH1	1:B:246:LEU:HD22	2.33	0.44
1:A:40:GLN:OE1	1:B:40[A]:GLN:NE2	2.51	0.44
1:B:103:ARG:HD2	4:B:747:HOH:O	2.17	0.43
2:D:137:SER:O	2:D:140:ILE:HB	2.18	0.43
1:B:69:ARG:HB3	1:B:81:ASP:HB2	1.99	0.43
2:D:9:HIS:HD2	2:D:11:PHE:N	2.13	0.43
2:D:77:MET:HE1	2:D:144:ALA:HA	1.99	0.43
2:C:33:LYS:HE3	2:C:124:GLN:NE2	2.33	0.43
1:A:88:ARG:NH2	1:B:92:GLU:OE2	2.53	0.42
1:A:88:ARG:NH2	1:B:88[A]:ARG:HD3	2.30	0.42
1:A:313:ALA:HA	1:A:318:PHE:CD2	2.54	0.42
2:D:65:VAL:CG1	2:D:136:LEU:HD21	2.49	0.42
2:D:77:MET:HE3	2:D:107:LEU:CD1	2.36	0.42
2:C:39:MET:CE	2:C:42:LYS:HE2	2.50	0.42
1:B:203:HIS:HE1	4:B:734:HOH:O	2.03	0.41
1:A:336:VAL:HG11	1:A:341:MET:HE3	2.02	0.41
1:B:145:GLN:H	1:B:145:GLN:HG3	1.52	0.41
2:C:39:MET:HE1	2:C:42:LYS:HE2	2.01	0.41
2:C:64:ARG:HD3	4:C:217:HOH:O	2.20	0.41
1:A:71:LYS:HE3	1:A:287:GLU:HB2	2.02	0.41
1:B:88[B]:ARG:HG2	4:B:761:HOH:O	2.21	0.41
2:C:110:GLN:HG3	4:C:247:HOH:O	2.20	0.41
1:A:26:ALA:O	1:A:222:LYS:HE2	2.20	0.41
1:A:375:SER:HB2	4:A:702:HOH:O	2.21	0.41
2:C:49:ASP:HA	2:C:52:LYS:HD2	2.03	0.41
1:A:341:MET:HE3	1:A:341:MET:HB2	1.86	0.41
1:A:122:ASN:HD22	1:A:122:ASN:HA	1.71	0.40
2:D:39[A]:MET:HE3	2:D:39[A]:MET:HB3	1.85	0.40
2:D:124:GLN:CD	2:D:124:GLN:H	2.29	0.40
1:A:9:PHE:CE1	1:A:378:PRO:HB2	2.56	0.40
1:A:341:MET:HG2	1:A:372:LEU:HD21	2.03	0.40
1:B:371:THR:HG22	1:B:372:LEU:N	2.35	0.40
1:A:198:ALA:HB3	3:A:500:PLP:C2	2.51	0.40
1:B:330:ALA:HB1	1:B:371:THR:HG21	2.03	0.40
2:C:138:GLU:OE1	2:C:138:GLU:HA	2.20	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	389/404 (96%)	379 (97%)	10 (3%)	0	100	100
1	B	403/404 (100%)	389 (96%)	13 (3%)	1 (0%)	43	42
2	C	146/150 (97%)	142 (97%)	4 (3%)	0	100	100
2	D	134/150 (89%)	131 (98%)	3 (2%)	0	100	100
All	All	1072/1108 (97%)	1041 (97%)	30 (3%)	1 (0%)	48	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	19	ALA

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	307/317 (97%)	296 (96%)	11 (4%)	31	31
1	B	319/317 (101%)	309 (97%)	10 (3%)	35	37
2	C	115/116 (99%)	110 (96%)	5 (4%)	26	25
2	D	106/116 (91%)	94 (89%)	12 (11%)	5	3
All	All	847/866 (98%)	809 (96%)	38 (4%)	26	23

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	GLN
1	A	80	ASP
1	A	115	SER
1	A	148	PRO
1	A	252	VAL
1	A	295	ASN
1	A	307	LEU
1	A	339	SER
1	A	341	MET
1	A	345	LEU
1	A	398	LEU
1	B	18	ASP
1	B	80	ASP
1	B	103	ARG
1	B	115	SER
1	B	145	GLN
1	B	148	PRO
1	B	237[A]	GLU
1	B	237[B]	GLU
1	B	246	LEU
1	B	365	GLU
2	C	61	CYS
2	C	71	VAL
2	C	102	LYS
2	C	146	GLN
2	C	147	VAL
2	D	18	GLU
2	D	36	GLN
2	D	39[A]	MET
2	D	39[B]	MET
2	D	40	LEU
2	D	51	LEU
2	D	58	ILE
2	D	77	MET
2	D	89	ARG
2	D	96	LEU
2	D	129	ARG
2	D	141	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	77	ASN
1	A	95	ASN
1	A	122	ASN
1	A	210	GLN
1	A	253	HIS
1	A	295	ASN
1	A	337	HIS
1	A	338	HIS
1	B	8	GLN
1	B	41	GLN
1	B	95	ASN
1	B	122	ASN
1	B	337	HIS
1	B	357	HIS
1	B	390	ASN
2	C	3	ASN
2	C	9	HIS
2	C	43	GLN
2	C	54	GLN
2	C	124	GLN
2	D	9	HIS
2	D	36	GLN
2	D	78	HIS
2	D	131	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 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	$\# Z > 2$
3	PLP	B	500	1	15,15,16	2.05	4 (26%)	21,22,23	1.76	6 (28%)
3	PLP	A	500	1	15,15,16	2.11	3 (20%)	21,22,23	2.19	9 (42%)

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	PLP	B	500	1	-	0/6/6/8	0/1/1/1
3	PLP	A	500	1	-	0/6/6/8	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	PLP	C4A-C4	-5.40	1.40	1.51
3	B	500	PLP	C4A-C4	-5.11	1.41	1.51
3	B	500	PLP	C2A-C2	-4.09	1.43	1.50
3	A	500	PLP	C2A-C2	-3.94	1.44	1.50
3	A	500	PLP	C2-N1	2.79	1.38	1.33
3	B	500	PLP	C6-N1	2.74	1.40	1.34
3	B	500	PLP	C5A-C5	-2.72	1.43	1.50

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	PLP	O4P-C5A-C5	4.26	117.35	109.36
3	A	500	PLP	C2A-C2-C3	3.88	125.33	120.80
3	A	500	PLP	C4A-C4-C5	-3.70	117.13	120.94
3	A	500	PLP	C3-C2-N1	-3.68	116.32	120.96
3	B	500	PLP	O4P-C5A-C5	3.23	115.42	109.36
3	B	500	PLP	O4P-P-O1P	3.20	115.08	106.44
3	B	500	PLP	C3-C2-N1	-3.18	116.94	120.96
3	A	500	PLP	C3-C4-C5	2.76	121.90	118.59

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	500	PLP	C4A-C4-C5	-2.62	118.24	120.94
3	A	500	PLP	O3-C3-C2	2.53	122.83	117.58
3	B	500	PLP	O2P-P-O4P	-2.48	100.20	106.67
3	B	500	PLP	O3P-P-O1P	2.41	120.24	110.83
3	A	500	PLP	O4P-P-O1P	2.27	112.57	106.44
3	A	500	PLP	O3P-P-O1P	2.27	119.67	110.83
3	A	500	PLP	C6-N1-C2	2.26	123.30	119.20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	500	PLP	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	392/404 (97%)	-0.50	8 (2%) 65 64	10, 18, 31, 65	1 (0%)
1	B	398/404 (98%)	-0.67	3 (0%) 82 82	8, 15, 29, 48	7 (1%)
2	C	145/150 (96%)	0.17	4 (2%) 55 54	11, 27, 50, 62	3 (2%)
2	D	137/150 (91%)	1.52	38 (27%) 1 1	24, 55, 75, 81	1 (0%)
All	All	1072/1108 (96%)	-0.22	53 (4%) 35 34	8, 19, 59, 81	12 (1%)

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	61	CYS	5.3
1	B	19	ALA	4.8
2	C	59	ALA	4.1
2	D	51	LEU	3.8
2	D	113	LEU	3.8
2	D	59	ALA	3.7
2	D	72	ALA	3.7
2	D	104	ALA	3.6
2	D	136	LEU	3.5
2	D	112	PRO	3.3
2	D	144	ALA	3.3
2	D	60	GLY	3.2
2	D	111	SER	3.1
1	A	55	PHE	3.1
2	D	115	LEU	3.1
1	A	56	ALA	3.0
2	D	140	ILE	3.0
2	D	114	ALA	3.0
2	D	69	TYR	3.0
2	D	145	LYS	2.8
2	D	48	PRO	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
2	C	147	VAL	2.8
2	D	109	ALA	2.7
2	D	58	ILE	2.7
2	D	12	GLY	2.5
1	A	400	VAL	2.5
2	D	17	ALA	2.5
2	D	47	LEU	2.5
2	D	70	THR	2.5
2	D	16	THR	2.4
1	A	48	GLY	2.4
2	D	55	ALA	2.4
1	B	18	ASP	2.4
2	D	39[A]	MET	2.4
2	D	127	ALA	2.3
2	D	67	LEU	2.3
1	A	88	ARG	2.3
2	D	13	THR	2.2
1	A	47	ALA	2.2
1	B	17	GLN	2.2
2	D	53	ALA	2.2
2	D	143	ALA	2.2
1	A	3	VAL	2.1
2	D	105	ALA	2.1
1	A	10	ARG	2.1
2	C	134[A]	ASN	2.1
2	D	14	THR	2.1
2	D	101	GLY	2.1
2	D	141	ILE	2.1
2	D	66	TRP	2.1
2	D	15	VAL	2.1
2	D	129	ARG	2.0
2	D	71	VAL	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	PLP	A	500	15/16	0.99	0.04	11,15,18,21	0
3	PLP	B	500	15/16	0.99	0.04	8,12,14,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.