



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 06:01 am GMT

PDB ID : 1GSS
Title : THREE-DIMENSIONAL STRUCTURE OF CLASS PI GLUTATHIONE S-TRANSFERASE FROM HUMAN PLACENTA IN COMPLEX WITH S-HEXYLGLUTATHIONE AT 2.8 ANGSTROMS RESOLUTION
Authors : Reinemer, P.; Dirr, H.W.; Ladenstein, R.; Lobello, M.; Federici, G.; Huber, R.; Parker, M.W.
Deposited on : 1992-05-28
Resolution : 2.80 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

<http://wwpdb.org/validation/2016/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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

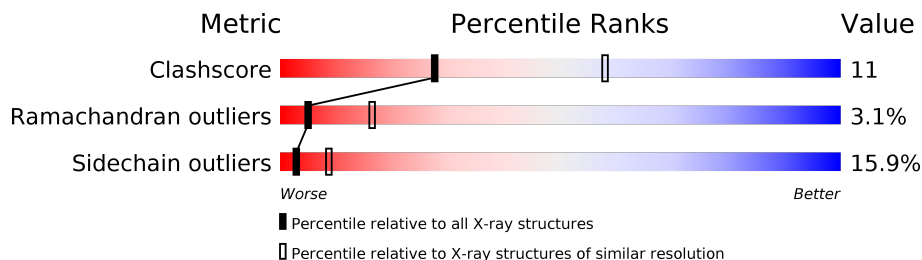
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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))
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)

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. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	209	
1	B	209	

2 Entry composition [i](#)

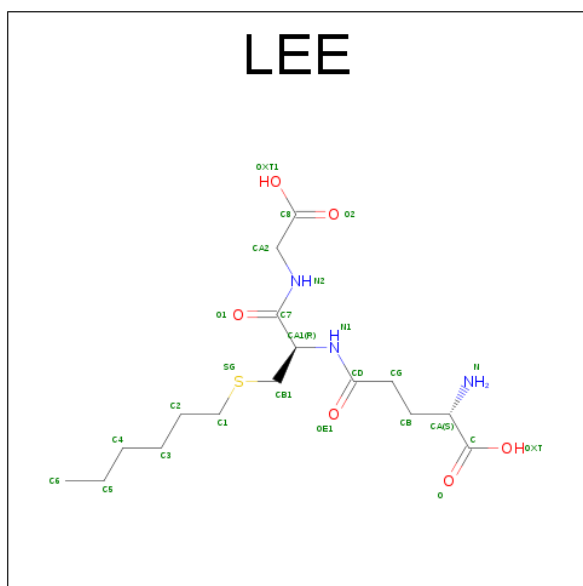
There are 3 unique types of molecules in this entry. The entry contains 3533 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GLUTATHIONE S-TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	209	Total	C	N	O	S	0	0	0
			1638	1052	272	308	6			
1	B	209	Total	C	N	O	S	0	0	0
			1638	1052	272	308	6			

- Molecule 2 is L-GAMMA-GLUTAMYL-S-HEXYL-L-CYSTEINYLGLYCINE (three-letter code: LEE) (formula: C₁₆H₂₉N₃O₆S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			26	16	3	6	1		
2	B	1	Total	C	N	O	S	0	0
			26	16	3	6	1		

- Molecule 3 is water.

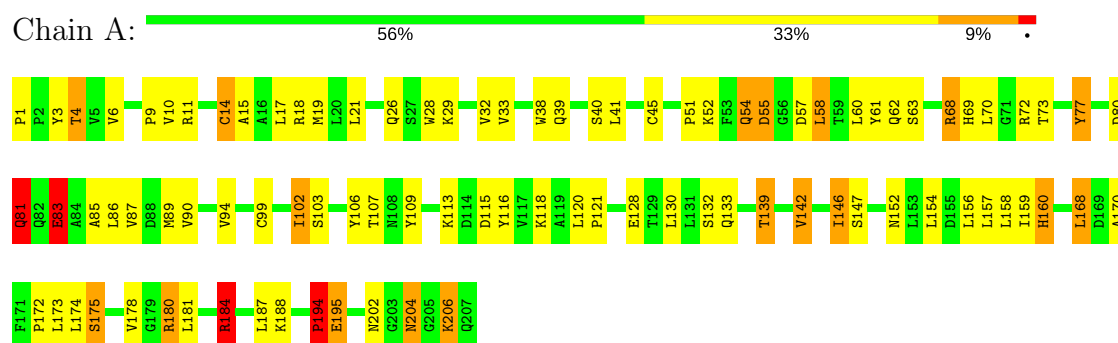
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	102	Total 102	O 102	0	0
3	B	103	Total 103	O 103	0	0

3 Residue-property plots

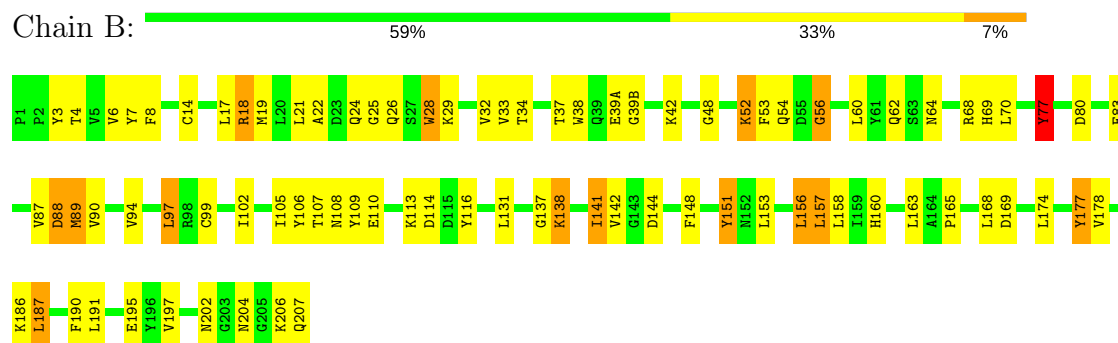
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

Note EDS was not executed.

• Molecule 1: GLUTATHIONE S-TRANSFERASE



• Molecule 1: GLUTATHIONE S-TRANSFERASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	60.50Å 60.50Å 238.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.80	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	EREF, X-PLOR	Depositor
R, R_{free}	0.196 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3533	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.83	1/1673 (0.1%)	1.67	24/2269 (1.1%)
1	B	0.83	0/1673	1.68	24/2269 (1.1%)
All	All	0.83	1/3346 (0.0%)	1.67	48/4538 (1.1%)

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	A	0	1
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	195	GLU	CB-CG	5.62	1.62	1.52

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	28	TRP	CD1-CG-CD2	9.04	113.53	106.30
1	B	89	MET	CA-CB-CG	-8.10	99.53	113.30
1	B	28	TRP	CE2-CD2-CG	-8.01	100.89	107.30
1	B	68	ARG	NE-CZ-NH1	7.95	124.27	120.30
1	A	38	TRP	CD1-CG-CD2	7.93	112.65	106.30
1	A	106	TYR	CB-CG-CD2	-7.65	116.41	121.00
1	A	77	TYR	CB-CG-CD2	-7.47	116.52	121.00
1	A	14	CYS	CA-CB-SG	-7.45	100.58	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	56	GLY	N-CA-C	-7.26	94.95	113.10
1	A	38	TRP	CE2-CD2-CG	-7.26	101.49	107.30
1	A	18	ARG	NE-CZ-NH2	-7.04	116.78	120.30
1	B	177	TYR	CB-CG-CD2	-7.02	116.79	121.00
1	A	195	GLU	CA-CB-CG	6.76	128.27	113.40
1	B	18	ARG	NE-CZ-NH2	-6.75	116.93	120.30
1	A	10	VAL	CG1-CB-CG2	-6.72	100.14	110.90
1	B	3	TYR	CB-CG-CD1	-6.70	116.98	121.00
1	A	28	TRP	CD1-CG-CD2	6.60	111.58	106.30
1	A	184	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	B	28	TRP	CB-CG-CD1	-6.58	118.44	127.00
1	A	160	HIS	CA-CB-CG	-6.47	102.61	113.60
1	A	28	TRP	CE2-CD2-CG	-6.44	102.15	107.30
1	B	163	LEU	CB-CA-C	-6.44	97.96	110.20
1	B	138	LYS	CB-CG-CD	6.43	128.32	111.60
1	B	77	TYR	CB-CG-CD2	-6.36	117.18	121.00
1	A	11	ARG	NE-CZ-NH1	6.35	123.47	120.30
1	A	89	MET	CG-SD-CE	-6.33	90.07	100.20
1	B	18	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	A	81	GLN	CA-CB-CG	6.21	127.07	113.40
1	A	11	ARG	NE-CZ-NH2	-6.04	117.28	120.30
1	A	116	TYR	CB-CG-CD1	-5.99	117.40	121.00
1	A	83	GLU	CA-CB-CG	5.94	126.47	113.40
1	B	38	TRP	CE2-CD2-CG	-5.78	102.68	107.30
1	B	28	TRP	CG-CD1-NE1	-5.70	104.40	110.10
1	B	28	TRP	CG-CD2-CE3	5.70	139.03	133.90
1	B	151	TYR	CB-CG-CD2	-5.67	117.60	121.00
1	A	194	PRO	CA-CB-CG	-5.64	93.28	104.00
1	A	1	PRO	CA-N-CD	-5.61	103.64	111.50
1	B	197	VAL	CG1-CB-CG2	-5.60	101.94	110.90
1	B	39(A)	GLU	CA-CB-CG	-5.60	101.08	113.40
1	B	163	LEU	CA-CB-CG	5.60	128.18	115.30
1	A	180	ARG	CA-C-N	5.50	129.30	117.20
1	B	52	LYS	CA-CB-CG	5.27	125.00	113.40
1	A	133	GLN	CA-CB-CG	5.26	124.98	113.40
1	A	57	ASP	N-CA-CB	-5.20	101.24	110.60
1	A	58	LEU	CA-CB-CG	5.13	127.11	115.30
1	B	107	THR	CA-CB-CG2	5.09	119.53	112.40
1	B	107	THR	N-CA-CB	-5.08	100.65	110.30
1	B	80	ASP	CA-CB-CG	5.07	124.56	113.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	184	ARG	Peptide
1	B	77	TYR	Sidechain

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	1638	0	1643	43	0
1	B	1638	0	1643	34	0
2	A	26	0	27	1	0
2	B	26	0	27	0	0
3	A	102	0	0	3	0
3	B	103	0	0	2	0
All	All	3533	0	3340	75	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 11.

All (75) 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:105:ILE:HG21	1:B:202:ASN:HD22	1.57	0.68
1:A:87:VAL:HG12	1:A:142:VAL:HG11	1.77	0.66
1:B:14:CYS:O	1:B:18:ARG:HG3	1.95	0.66
1:A:4:THR:HG23	1:A:29:LYS:HB2	1.78	0.65
1:B:105:ILE:HG12	1:B:202:ASN:ND2	2.12	0.64
1:B:109:TYR:O	1:B:113:LYS:HB2	1.98	0.64
1:A:174:LEU:O	1:A:178:VAL:HG13	1.99	0.63
1:B:90:VAL:HG11	1:B:141:ILE:HG21	1.80	0.63
1:A:175:SER:O	1:A:178:VAL:HG22	2.01	0.61
1:B:97:LEU:HD12	1:B:156:LEU:HD21	1.85	0.58
1:B:105:ILE:HG21	1:B:202:ASN:ND2	2.20	0.57
1:A:157:LEU:HD22	1:A:168:LEU:HD11	1.87	0.56
1:B:153:LEU:HD23	1:B:177:TYR:CE2	2.41	0.56
1:A:21:LEU:HB3	1:A:26:GLN:HB2	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:VAL:HB	1:A:52:LYS:HB3	1.88	0.54
1:B:83:GLU:O	1:B:87:VAL:HG13	2.07	0.54
1:B:8:PHE:HA	1:B:33:VAL:O	2.08	0.54
1:A:6:VAL:HG21	1:A:52:LYS:HD3	1.92	0.52
1:A:85:ALA:HB2	1:B:69:HIS:HD2	1.73	0.52
1:B:90:VAL:HG11	1:B:141:ILE:CG2	2.39	0.52
1:A:181:LEU:O	1:A:187:LEU:HD23	2.12	0.50
1:A:85:ALA:HB1	1:B:60:LEU:HD21	1.93	0.50
1:A:45:CYS:HA	1:A:61:TYR:CZ	2.47	0.49
1:A:54:GLN:HA	1:A:58:LEU:O	2.12	0.49
1:A:146:ILE:HD12	1:A:184:ARG:CZ	2.43	0.49
1:B:204:ASN:ND2	1:B:206:LYS:HG2	2.28	0.48
1:A:154:LEU:O	1:A:158:LEU:HG	2.14	0.48
1:B:19:MET:HB3	1:B:187:LEU:HD21	1.94	0.47
1:A:14:CYS:SG	1:A:51:PRO:HB3	2.56	0.46
1:B:190:PHE:HA	3:B:253:HOH:O	2.16	0.46
1:A:94:VAL:HG13	1:A:156:LEU:HD22	1.98	0.46
1:B:32:VAL:HG22	3:B:225:HOH:O	2.15	0.46
1:A:69:HIS:HA	1:A:72:ARG:NH1	2.31	0.46
1:B:53:PHE:CE2	1:B:70:LEU:HD11	2.51	0.46
1:A:128:GLU:HG3	1:A:173:LEU:HD22	1.98	0.45
1:A:184:ARG:NH1	3:A:293:HOH:O	2.48	0.45
1:A:3:TYR:HE1	1:A:55:ASP:OD1	1.99	0.45
1:B:157:LEU:HD23	1:B:178:VAL:CG2	2.46	0.45
1:B:24:GLN:O	1:B:26:GLN:N	2.49	0.45
1:A:204:ASN:HD21	1:A:206:LYS:HB2	1.82	0.45
1:A:90:VAL:HG22	1:A:130:LEU:HD13	1.99	0.45
1:B:148:PHE:HA	1:B:151:TYR:HD2	1.82	0.45
1:B:116:TYR:HH	1:B:160:HIS:HE2	1.65	0.45
1:A:170:ALA:O	1:A:172:PRO:HD3	2.17	0.44
1:A:19:MET:HB3	1:A:187:LEU:HD11	1.99	0.44
1:B:87:VAL:HA	1:B:142:VAL:HG21	2.00	0.44
1:A:109:TYR:CG	1:A:206:LYS:HG2	2.53	0.44
1:B:6:VAL:HG22	1:B:52:LYS:HB3	1.99	0.44
1:B:22:ALA:HB2	1:B:28:TRP:HH2	1.82	0.44
1:A:87:VAL:HG13	1:A:142:VAL:HG21	2.00	0.44
1:B:94:VAL:HG22	1:B:153:LEU:HD12	2.00	0.44
1:A:181:LEU:O	1:A:184:ARG:HG2	2.17	0.43
1:A:15:ALA:O	1:A:19:MET:HG3	2.18	0.43
1:B:22:ALA:HB2	1:B:28:TRP:CH2	2.53	0.43
1:A:202:ASN:HB2	1:A:206:LYS:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:VAL:HG13	1:A:142:VAL:O	2.18	0.43
1:A:103:SER:O	1:A:107:THR:HB	2.19	0.43
1:A:9:PRO:HD3	1:A:33:VAL:O	2.19	0.43
1:B:4:THR:HB	1:B:54:GLN:HB2	2.00	0.42
1:A:139:THR:O	1:A:180:ARG:NH1	2.53	0.42
1:B:64:ASN:HA	1:B:64:ASN:HD22	1.52	0.42
1:A:9:PRO:HG3	1:A:32:VAL:CG1	2.49	0.42
2:A:208:LEE:HB31	3:A:263:HOH:O	2.19	0.42
1:A:81:GLN:HB3	1:A:81:GLN:HE21	1.71	0.42
1:A:188:LYS:HB3	1:A:188:LYS:HE3	1.83	0.42
1:A:86:LEU:O	1:A:90:VAL:HG23	2.20	0.41
1:A:120:LEU:HD11	1:A:160:HIS:HD2	1.85	0.41
1:A:68:ARG:HD2	1:B:88:ASP:OD2	2.19	0.41
1:B:131:LEU:HD12	1:B:137:GLY:HA2	2.01	0.41
1:A:80:ASP:OD1	1:A:83:GLU:HB2	2.21	0.41
1:B:105:ILE:HG22	1:B:106:TYR:CD2	2.56	0.41
1:B:202:ASN:N	1:B:202:ASN:OD1	2.55	0.40
1:B:42:LYS:HG3	1:B:48:GLY:O	2.22	0.40
1:A:156:LEU:O	1:A:159:ILE:HG12	2.21	0.40
1:A:102:ILE:HD12	3:A:310:HOH:O	2.22	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	207/209 (99%)	186 (90%)	15 (7%)	6 (3%)	5	18
1	B	207/209 (99%)	179 (86%)	21 (10%)	7 (3%)	4	15
All	All	414/418 (99%)	365 (88%)	36 (9%)	13 (3%)	5	16

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	25	GLY
1	A	142	VAL
1	B	108	ASN
1	A	195	GLU
1	B	114	ASP
1	B	165	PRO
1	A	62	GLN
1	A	204	ASN
1	A	63	SER
1	B	39(B)	GLY
1	B	62	GLN
1	A	194	PRO
1	B	56	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	176/176 (100%)	147 (84%)	29 (16%)	2	8
1	B	176/176 (100%)	149 (85%)	27 (15%)	3	9
All	All	352/352 (100%)	296 (84%)	56 (16%)	3	9

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	17	LEU
1	A	39	GLN
1	A	40	SER
1	A	41	LEU
1	A	54	GLN
1	A	55	ASP
1	A	60	LEU
1	A	68	ARG
1	A	70	LEU
1	A	73	THR

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Mol	Chain	Res	Type
1	A	77	TYR
1	A	81	GLN
1	A	83	GLU
1	A	99	CYS
1	A	102	ILE
1	A	113	LYS
1	A	115	ASP
1	A	118	LYS
1	A	121	PRO
1	A	132	SER
1	A	139	THR
1	A	146	ILE
1	A	147	SER
1	A	152	ASN
1	A	168	LEU
1	A	175	SER
1	A	194	PRO
1	A	206	LYS
1	B	7	TYR
1	B	17	LEU
1	B	21	LEU
1	B	29	LYS
1	B	34	THR
1	B	37	THR
1	B	77	TYR
1	B	88	ASP
1	B	89	MET
1	B	97	LEU
1	B	99	CYS
1	B	102	ILE
1	B	110	GLU
1	B	138	LYS
1	B	141	ILE
1	B	144	ASP
1	B	156	LEU
1	B	157	LEU
1	B	158	LEU
1	B	168	LEU
1	B	169	ASP
1	B	174	LEU
1	B	186	LYS
1	B	187	LEU

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Mol	Chain	Res	Type
1	B	191	LEU
1	B	195	GLU
1	B	207	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	81	GLN
1	B	26	GLN
1	B	64	ASN
1	B	69	HIS
1	B	82	GLN
1	B	91	ASN
1	B	152	ASN
1	B	202	ASN

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](#)

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
2	LEE	A	208	-	17,25,25	0.68	0	20,30,30	3.44	9 (45%)
2	LEE	B	208	-	17,25,25	1.67	1 (5%)	20,30,30	1.84	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	LEE	A	208	-	-	0/24/30/30	0/0/0/0
2	LEE	B	208	-	-	0/24/30/30	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	208	LEE	OE1-CD	-6.27	1.10	1.23

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	208	LEE	CG-CD-N1	-7.18	103.19	115.82
2	A	208	LEE	CA1-CB1-SG	-5.04	100.28	113.37
2	A	208	LEE	O1-C7-N2	-5.00	113.45	123.07
2	A	208	LEE	CB1-CA1-N1	-4.61	99.04	110.92
2	B	208	LEE	CG-CD-N1	-4.45	107.99	115.82
2	B	208	LEE	CB1-SG-C1	-3.90	90.48	102.29
2	A	208	LEE	CB-CG-CD	-3.47	105.34	113.18
2	B	208	LEE	CA2-N2-C7	-3.22	118.03	122.39
2	B	208	LEE	CB-CG-CD	-2.07	108.50	113.18
2	A	208	LEE	C7-CA1-N1	-2.04	105.58	111.20
2	A	208	LEE	CB1-CA1-C7	4.35	119.56	109.55
2	A	208	LEE	CA1-C7-N2	5.81	128.37	116.51
2	A	208	LEE	OE1-CD-N1	6.03	133.38	122.97

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
2	A	208	LEE	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.