



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:02 AM GMT

PDB ID : 3GZC
Title : Structure of human selenocysteine lyase
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Deposited on : 2009-04-07
Resolution : 2.10 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

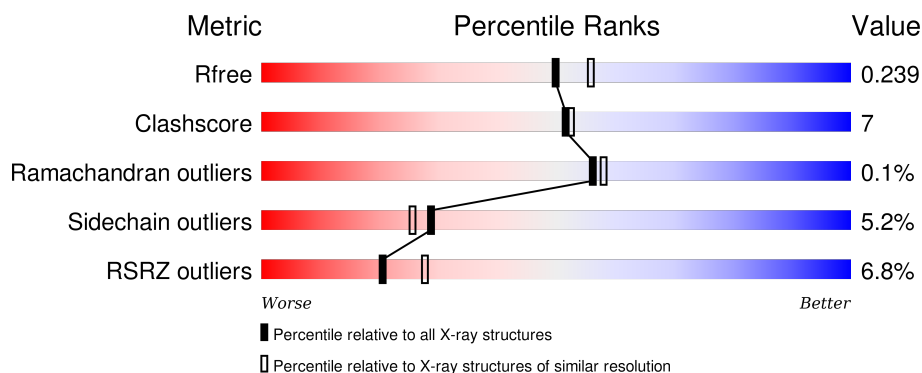
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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.

Mol	Chain	Length	Quality of chain
1	A	440	<div> <div>5%</div> <div>78% 12% • 8%</div> </div>
1	B	440	<div> <div>8%</div> <div>76% 13% • 9%</div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6684 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.

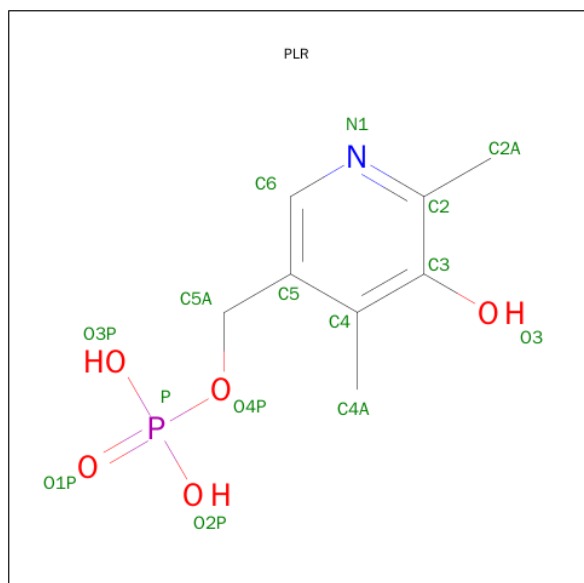
- Molecule 1 is a protein called Selenocysteine lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	403	Total	C	N	O	S	0	0	0
			3103	1948	558	583	14			
1	B	402	Total	C	N	O	S	0	0	0
			3094	1943	556	581	14			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	6	SER	-	EXPRESSION TAG	UNP Q96I15
A	7	MET	-	EXPRESSION TAG	UNP Q96I15
A	175	THR	ALA	SEE REMARK 999	UNP Q96I15
B	6	SER	-	EXPRESSION TAG	UNP Q96I15
B	7	MET	-	EXPRESSION TAG	UNP Q96I15
B	175	THR	ALA	SEE REMARK 999	UNP Q96I15

- Molecule 2 is (5-HYDROXY-4,6-DIMETHYLPYRIDIN-3-YL)METHYL DIHYDROGEN PHOSPHATE (three-letter code: PLR) (formula: C₈H₁₂NO₅P).



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

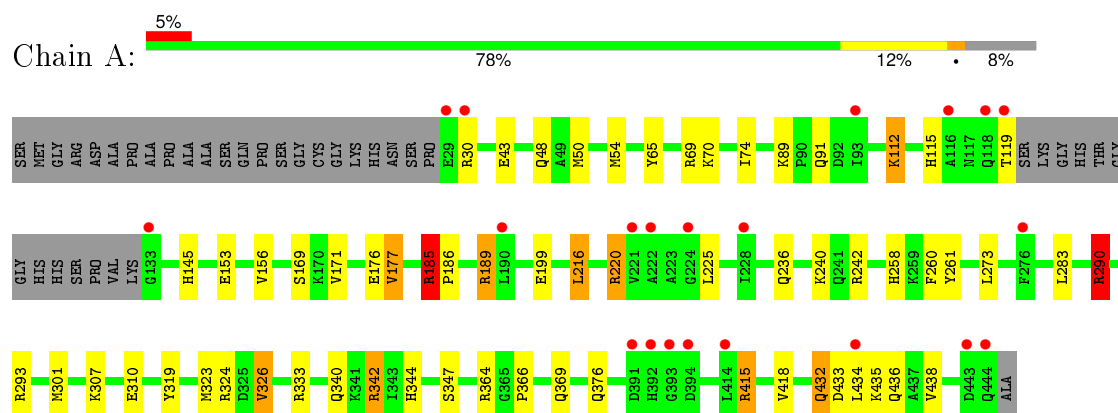
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	269	Total	O	0	0
			269	269		
3	B	188	Total	O	0	0
			188	188		

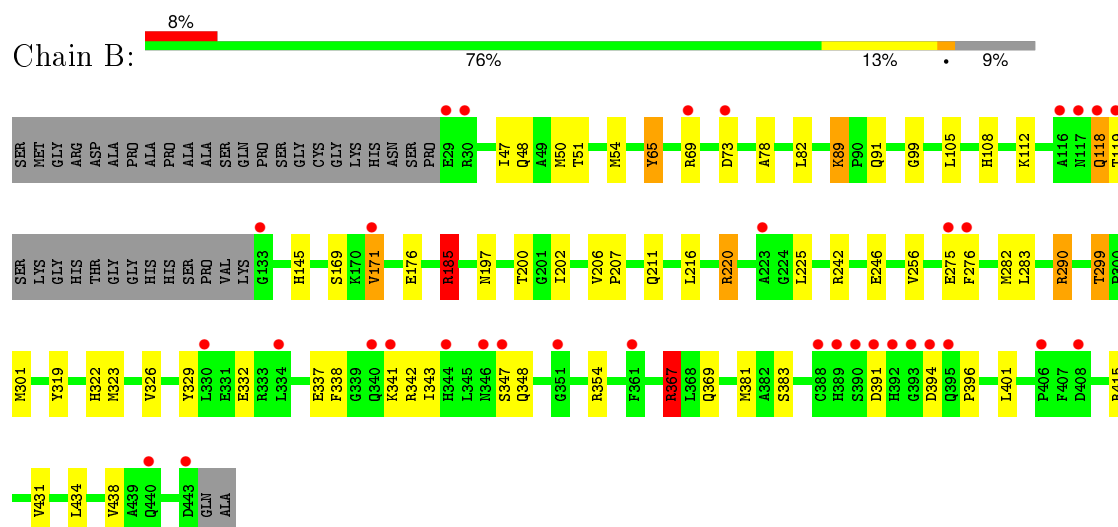
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Selenocysteine lyase



• Molecule 1: Selenocysteine lyase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.22Å 85.81Å 188.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.73 – 2.10 19.72 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.73-2.10) 99.9 (19.72-2.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.86 (at 2.09Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.179 , 0.235 0.184 , 0.239	Depositor DCC
R_{free} test set	2844 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 47.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Outliers	0 of 56897 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6684	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 3.63% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: 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	0.92	3/3164 (0.1%)	0.92	11/4295 (0.3%)
1	B	0.79	0/3155	0.84	6/4283 (0.1%)
All	All	0.86	3/6319 (0.0%)	0.88	17/8578 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	326	VAL	CB-CG2	-6.21	1.39	1.52
1	A	260	PHE	CE2-CZ	5.33	1.47	1.37
1	A	310	GLU	CG-CD	5.02	1.59	1.51

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	189	ARG	NE-CZ-NH1	9.50	125.05	120.30
1	A	189	ARG	NE-CZ-NH2	-9.26	115.67	120.30
1	A	220	ARG	NE-CZ-NH2	-8.89	115.85	120.30
1	A	324	ARG	NE-CZ-NH2	8.87	124.74	120.30
1	A	220	ARG	NE-CZ-NH1	8.68	124.64	120.30
1	A	324	ARG	NE-CZ-NH1	-8.01	116.29	120.30
1	B	220	ARG	NE-CZ-NH2	-7.86	116.37	120.30
1	A	415	ARG	NE-CZ-NH1	7.73	124.16	120.30
1	B	220	ARG	NE-CZ-NH1	7.70	124.15	120.30
1	B	367	ARG	NE-CZ-NH1	6.91	123.76	120.30
1	B	367	ARG	CG-CD-NE	6.00	124.40	111.80
1	B	185	ARG	NE-CZ-NH2	5.91	123.25	120.30
1	A	185	ARG	NE-CZ-NH2	5.62	123.11	120.30
1	A	333	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	B	367	ARG	NE-CZ-NH2	-5.53	117.54	120.30
1	A	185	ARG	NE-CZ-NH1	-5.37	117.61	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	290	ARG	NE-CZ-NH1	5.01	122.81	120.30

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	3103	0	3099	41	0
1	B	3094	0	3090	47	0
2	A	15	0	7	1	0
2	B	15	0	8	1	0
3	A	269	0	0	15	0
3	B	188	0	0	8	0
All	All	6684	0	6204	85	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 7.

All (85) 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:50:MET:HE1	1:A:301:MET:CE	1.70	1.19
1:B:50:MET:CE	1:B:301:MET:HE1	1.86	1.05
1:A:50:MET:CE	1:A:301:MET:HE1	1.89	1.02
1:B:50:MET:HE1	1:B:301:MET:CE	1.91	0.99
1:B:50:MET:HE1	1:B:301:MET:HE1	0.98	0.96
1:A:50:MET:HE1	1:A:301:MET:HE1	0.94	0.93
1:B:290:ARG:HH11	1:B:290:ARG:HG2	1.45	0.82
1:A:50:MET:CE	1:A:301:MET:CE	2.51	0.82
1:B:415:ARG:HD2	3:B:809:HOH:O	1.85	0.75
1:B:367:ARG:HG3	1:B:367:ARG:HH11	1.55	0.72
1:B:47:ILE:O	1:B:51:THR:HG23	1.93	0.69
1:B:169:SER:OG	1:B:171:VAL:HG13	1.96	0.66
1:A:74:ILE:HD13	3:A:605:HOH:O	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:337:GLU:HG3	1:B:431:VAL:HG11	1.80	0.62
1:A:319:TYR:HB3	1:A:323:MET:CE	2.31	0.60
1:A:344:HIS:NE2	1:A:364:ARG:HD2	2.17	0.60
1:A:115:HIS:HE1	3:A:639:HOH:O	1.87	0.58
1:B:348:GLN:HE21	1:B:354:ARG:NH2	2.03	0.56
1:B:197:ASN:HD22	1:B:200:THR:H	1.53	0.56
1:A:145:HIS:HB2	2:A:500:PLR:H2A3	1.87	0.56
1:A:50:MET:CE	1:A:301:MET:HE2	2.35	0.55
1:B:242:ARG:HD3	3:B:786:HOH:O	2.06	0.55
1:A:70:LYS:HE2	3:A:842:HOH:O	2.05	0.55
1:A:366:PRO:O	1:A:369:GLN:NE2	2.39	0.54
1:B:202:ILE:HD11	1:B:347:SER:O	2.07	0.53
1:A:91:GLN:HB2	3:A:457:HOH:O	2.09	0.53
1:A:242:ARG:NE	3:A:744:HOH:O	2.43	0.52
1:B:369:GLN:HG2	1:B:394:ASP:OD2	2.11	0.51
1:B:50:MET:CE	1:B:301:MET:CE	2.70	0.51
1:B:299:THR:CG2	3:B:3:HOH:O	2.59	0.51
1:A:153:GLU:O	1:A:156:VAL:HG12	2.11	0.51
1:B:319:TYR:C	1:B:323:MET:HE2	2.30	0.50
1:B:319:TYR:O	1:B:323:MET:HE2	2.13	0.49
1:B:78:ALA:O	1:B:82:LEU:HD23	2.12	0.49
1:A:347:SER:HB2	3:A:628:HOH:O	2.13	0.48
1:B:299:THR:HG23	3:B:3:HOH:O	2.12	0.48
1:A:415:ARG:HD2	3:A:801:HOH:O	2.13	0.48
1:A:30:ARG:HD3	3:A:767:HOH:O	2.13	0.47
1:B:434:LEU:O	1:B:438:VAL:HG13	2.14	0.47
1:B:108:HIS:CD2	1:B:282:MET:HG2	2.50	0.47
1:B:396:PRO:HG2	1:B:401:LEU:HD11	1.97	0.47
1:B:337:GLU:HG3	1:B:431:VAL:CG1	2.43	0.47
1:A:216:LEU:HD22	1:A:220:ARG:HG3	1.97	0.47
1:B:176:GLU:HG3	3:B:621:HOH:O	2.14	0.46
1:A:43:GLU:HG2	1:A:261:TYR:HB3	1.97	0.46
1:B:50:MET:HE3	1:B:301:MET:HA	1.98	0.46
1:B:322:HIS:O	1:B:326:VAL:HG13	2.14	0.46
1:A:186:PRO:O	1:A:189:ARG:NH2	2.48	0.46
1:A:91:GLN:CB	3:A:457:HOH:O	2.63	0.46
1:B:145:HIS:HB2	2:B:500:PLR:H2A3	1.98	0.45
1:A:112:LYS:HA	1:A:112:LYS:HD2	1.88	0.45
1:B:118:GLN:O	1:B:119:THR:OG1	2.27	0.45
1:B:391:ASP:OD2	1:B:391:ASP:N	2.50	0.45
1:A:185:ARG:O	1:A:220:ARG:NH2	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:273:LEU:HD23	1:A:290:ARG:NH2	2.32	0.45
1:A:65:TYR:O	1:A:69:ARG:HG2	2.17	0.45
1:B:89:LYS:NZ	3:B:572:HOH:O	2.44	0.44
1:B:383:SER:HB2	3:B:599:HOH:O	2.17	0.44
1:A:169:SER:OG	1:A:171:VAL:HG12	2.17	0.44
1:B:99:GLY:HA3	1:B:256:VAL:HG22	1.98	0.44
1:B:112:LYS:NZ	3:B:832:HOH:O	2.43	0.44
1:A:432:GLN:NE2	1:A:435:LYS:HD3	2.31	0.44
1:B:329:TYR:HA	1:B:332:GLU:HG2	1.99	0.44
1:A:258:HIS:HD2	3:A:453:HOH:O	2.00	0.44
1:A:156:VAL:HG11	3:A:586:HOH:O	2.17	0.44
1:A:290:ARG:HG2	3:A:758:HOH:O	2.18	0.44
1:A:342:ARG:NH2	3:A:754:HOH:O	2.51	0.43
1:B:185:ARG:O	1:B:220:ARG:NH2	2.51	0.43
1:B:108:HIS:CG	1:B:282:MET:HG2	2.54	0.43
1:B:338:PHE:HB2	1:B:343:ILE:HD11	1.99	0.43
1:A:434:LEU:O	1:A:438:VAL:HG13	2.19	0.42
1:B:50:MET:CE	1:B:301:MET:HA	2.49	0.42
1:B:275:GLU:HG3	1:B:276:PHE:CD1	2.55	0.42
1:A:236:GLN:O	1:A:240:LYS:HD2	2.19	0.42
1:A:376:GLN:HG2	3:A:601:HOH:O	2.19	0.41
1:B:206:VAL:N	1:B:207:PRO:CD	2.83	0.41
1:B:290:ARG:HH11	1:B:290:ARG:CG	2.22	0.41
1:B:89:LYS:HB3	1:B:91:GLN:OE1	2.20	0.41
1:B:105:LEU:HD13	1:B:282:MET:HG3	2.02	0.41
1:A:65:TYR:CZ	1:B:381:MET:HA	2.56	0.41
1:A:326:VAL:CG2	1:A:418:VAL:HG21	2.51	0.41
1:A:433:ASP:O	1:A:436:GLN:HG3	2.21	0.40
1:A:54:MET:HE1	1:B:54:MET:HE3	2.02	0.40
1:A:54:MET:HE3	1:B:54:MET:HE1	2.04	0.40
1:A:177:VAL:HG13	3:A:464:HOH:O	2.21	0.40

There are no symmetry-related clashes.

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	A	399/440 (91%)	390 (98%)	9 (2%)	0	100	100
1	B	398/440 (90%)	388 (98%)	9 (2%)	1 (0%)	46	45
All	All	797/880 (91%)	778 (98%)	18 (2%)	1 (0%)	56	58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	65	TYR

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	334/360 (93%)	317 (95%)	17 (5%)	29	26
1	B	333/360 (92%)	315 (95%)	18 (5%)	27	24
All	All	667/720 (93%)	632 (95%)	35 (5%)	29	25

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	GLN
1	A	89	LYS
1	A	112	LYS
1	A	119	THR
1	A	176	GLU
1	A	177	VAL
1	A	185	ARG
1	A	199	GLU
1	A	216	LEU
1	A	225	LEU
1	A	283	LEU
1	A	290	ARG

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Mol	Chain	Res	Type
1	A	293	ARG
1	A	307	LYS
1	A	340	GLN
1	A	342	ARG
1	A	432	GLN
1	B	48	GLN
1	B	65	TYR
1	B	69	ARG
1	B	73	ASP
1	B	89	LYS
1	B	118	GLN
1	B	171	VAL
1	B	185	ARG
1	B	211	GLN
1	B	216	LEU
1	B	225	LEU
1	B	246	GLU
1	B	283	LEU
1	B	290	ARG
1	B	299	THR
1	B	341	LYS
1	B	342	ARG
1	B	367	ARG

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	115	HIS
1	A	241	GLN
1	B	113	HIS
1	B	118	GLN
1	B	197	ASN
1	B	211	GLN
1	B	314	GLN
1	B	348	GLN
1	B	392	HIS

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 carbohydrates 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	$\# Z > 2$
2	PLR	A	500	1	15,15,15	0.76	0	21,22,22	1.66	8 (38%)
2	PLR	B	500	1	15,15,15	0.96	0	21,22,22	2.29	8 (38%)

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	A	500	1	-	0/6/6/6	0/1/1/1
2	PLR	B	500	1	-	0/6/6/6	0/1/1/1

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	PLR	C4A-C4-C3	-3.57	113.89	120.36
2	B	500	PLR	C5-C6-N1	-2.47	119.58	123.86
2	A	500	PLR	O3P-P-O4P	-2.39	99.69	106.56
2	B	500	PLR	C5A-C5-C6	-2.26	115.01	119.28
2	A	500	PLR	C3-C2-N1	-2.04	117.80	120.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	PLR	O3P-P-O1P	2.04	117.15	110.58
2	A	500	PLR	C5A-C5-C4	2.21	124.58	121.65
2	A	500	PLR	O3-C3-C2	2.24	121.55	117.66
2	A	500	PLR	C6-N1-C2	2.29	123.95	119.28
2	B	500	PLR	O3-C3-C2	2.50	122.01	117.66
2	B	500	PLR	O4P-C5A-C5	2.72	113.49	108.99
2	A	500	PLR	O4P-C5A-C5	2.73	113.50	108.99
2	B	500	PLR	C6-N1-C2	2.74	124.87	119.28
2	A	500	PLR	C2A-C2-C3	3.33	125.06	121.04
2	B	500	PLR	C5A-C5-C4	3.40	126.16	121.65
2	B	500	PLR	C4A-C4-C5	6.36	127.51	120.88

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	PLR	1	0
2	B	500	PLR	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	403/440 (91%)	0.14	21 (5%) 31 39	25, 31, 44, 58	0
1	B	402/440 (91%)	0.32	34 (8%) 13 18	23, 32, 44, 61	0
All	All	805/880 (91%)	0.23	55 (6%) 20 28	23, 31, 44, 61	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	393	GLY	8.0
1	A	119	THR	7.1
1	B	29	GLU	5.5
1	A	276	PHE	4.6
1	A	443	ASP	4.4
1	B	351	GLY	4.3
1	B	118	GLN	4.1
1	B	389	HIS	4.1
1	B	69	ARG	4.1
1	B	171	VAL	4.0
1	A	444	GLN	3.8
1	A	29	GLU	3.8
1	A	394	ASP	3.7
1	A	118	GLN	3.6
1	B	347	SER	3.6
1	B	408	ASP	3.5
1	B	394	ASP	3.4
1	B	391	ASP	3.4
1	B	392	HIS	3.4
1	B	341	LYS	3.3
1	B	443	ASP	3.3
1	B	30	ARG	3.2
1	B	133	GLY	3.0
1	A	391	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	334	LEU	2.9
1	B	223	ALA	2.8
1	A	392	HIS	2.8
1	B	119	THR	2.8
1	B	340	GLN	2.7
1	B	117	ASN	2.6
1	A	133	GLY	2.6
1	B	330	LEU	2.5
1	A	30	ARG	2.5
1	A	393	GLY	2.4
1	B	344	HIS	2.4
1	A	116	ALA	2.4
1	A	222	ALA	2.4
1	A	414	LEU	2.4
1	B	406	PRO	2.3
1	A	93	ILE	2.3
1	B	276	PHE	2.3
1	B	116	ALA	2.3
1	A	434	LEU	2.3
1	B	361	PHE	2.2
1	B	73	ASP	2.2
1	A	228	ILE	2.2
1	B	388	CYS	2.1
1	B	346	ASN	2.1
1	A	221	VAL	2.1
1	B	395	GLN	2.1
1	B	440	GLN	2.1
1	A	190	LEU	2.0
1	A	224	GLY	2.0
1	B	390	SER	2.0
1	B	275	GLU	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
2	PLR	B	500	15/15	0.97	0.14	1.42	23,31,33,35	0
2	PLR	A	500	15/15	0.97	0.11	0.16	22,23,26,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.