



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:27 PM GMT

PDB ID : 4EZL
Title : Potent and Selective Inhibitors of PI3K-delta: Obtaining Isoform Selectivity from the Affinity Pocket and Tryptophan Shelf
Authors : Murray, J.M.
Deposited on : 2012-05-02
Resolution : 2.94 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

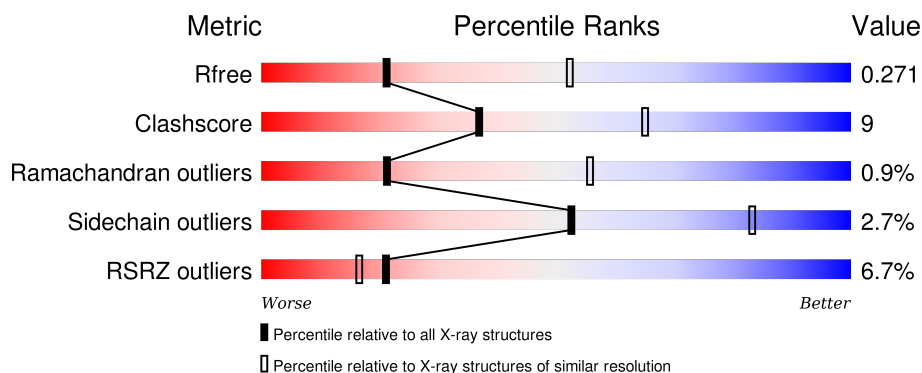
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.94 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	2057 (2.98-2.90)
Clashscore	102246	2308 (2.98-2.90)
Ramachandran outliers	100387	2245 (2.98-2.90)
Sidechain outliers	100360	2247 (2.98-2.90)
RSRZ outliers	91569	2065 (2.98-2.90)

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	966	<div> <div>6%</div> <div>66%</div> <div>18%</div> <div>•</div> <div>14%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6770 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.

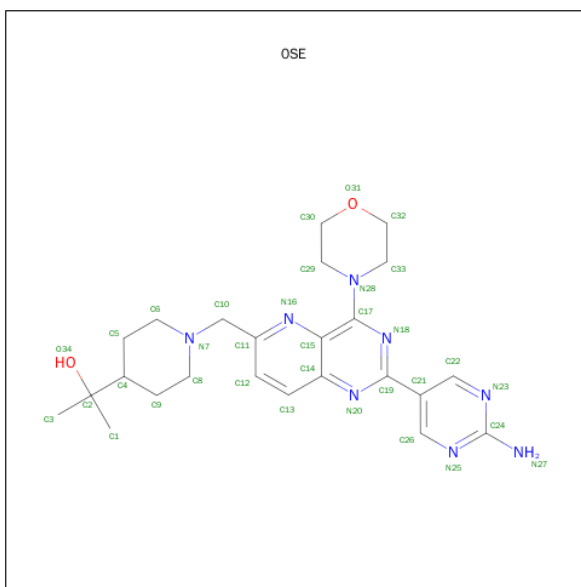
- Molecule 1 is a protein called Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic sub-unit gamma isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	834	Total	C	N	O	S	0	0	0
			6735	4324	1149	1228	34			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	143	MET	-	EXPRESSION TAG	UNP P48736
A	802	THR	LYS	ENGINEERED MUTATION	UNP P48736
A	1103	HIS	-	EXPRESSION TAG	UNP P48736
A	1104	HIS	-	EXPRESSION TAG	UNP P48736
A	1105	HIS	-	EXPRESSION TAG	UNP P48736
A	1106	HIS	-	EXPRESSION TAG	UNP P48736
A	1107	HIS	-	EXPRESSION TAG	UNP P48736
A	1108	HIS	-	EXPRESSION TAG	UNP P48736

- Molecule 2 is 2-(1-{[2-(2-AMINOPYRIMIDIN-5-YL)-4-(MORPHOLIN-4-YL)PYRIDO[3,2-D]PYRIMIDIN-6-YL]METHYL}PIPERIDIN-4-YL)PROPAN-2-OL (three-letter code: 0SE) (formula: C₂₄H₃₂N₈O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			34	24	8	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O	0	0
			1	1		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	141.33Å 67.19Å 106.56Å 90.00° 96.10° 90.00°	Depositor
Resolution (Å)	31.14 – 2.94 31.14 – 2.93	Depositor EDS
% Data completeness (in resolution range)	94.9 (31.14-2.94) 94.9 (31.14-2.93)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.95Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.7.2_869)	Depositor
R, R_{free}	0.244 , 0.271 0.242 , 0.271	Depositor DCC
R_{free} test set	1042 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	79.0	Xtriage
Anisotropy	0.066	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 56.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 20407 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6770	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.05% 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: 0SE

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	0.39	0/6878	0.77	9/9308 (0.1%)

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

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	832	PHE	CB-CG-CD2	-9.74	113.98	120.80
1	A	832	PHE	CB-CG-CD1	8.59	126.81	120.80
1	A	1042	LEU	CB-CG-CD2	-8.46	96.63	111.00
1	A	1092	LEU	CB-CA-C	-8.44	94.16	110.20
1	A	1092	LEU	N-CA-C	8.30	133.40	111.00
1	A	700	ARG	NE-CZ-NH1	-7.26	116.67	120.30
1	A	779	LEU	CA-CB-CG	6.62	130.52	115.30
1	A	700	ARG	NE-CZ-NH2	6.56	123.58	120.30
1	A	477	ARG	CA-CB-CG	5.32	125.10	113.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	899	THR	Peptide

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	6735	0	6755	127	0
2	A	34	0	32	3	0
3	A	1	0	0	0	0
All	All	6770	0	6787	127	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 9.

All (127) 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:1035:LEU:HD12	1:A:1048:ILE:HD13	1.49	0.94
1:A:1045:LYS:O	1:A:1049:GLU:HG3	1.69	0.91
1:A:1035:LEU:HB3	1:A:1042:LEU:CD2	2.06	0.85
1:A:564:LEU:HD11	1:A:1048:ILE:HG22	1.58	0.84
1:A:806:SER:O	1:A:807:LYS:CB	2.25	0.82
1:A:642:ALA:O	1:A:646:GLN:HG3	1.80	0.81
1:A:808:LYS:HG3	1:A:833:LYS:HE2	1.64	0.79
1:A:1035:LEU:HB3	1:A:1042:LEU:HD23	1.67	0.77
1:A:751:SER:O	1:A:752:LEU:HD22	1.89	0.73
1:A:149:ALA:O	1:A:153:GLN:HG3	1.90	0.72
1:A:1035:LEU:HB3	1:A:1042:LEU:HD21	1.73	0.71
1:A:382:PHE:CE2	1:A:398:ARG:HD3	2.28	0.69
1:A:1000:LYS:HG2	1:A:1000:LYS:O	1.91	0.69
1:A:745:VAL:HG12	1:A:811:LEU:HD11	1.76	0.68
1:A:564:LEU:HD21	1:A:1048:ILE:HG21	1.76	0.67
1:A:804:MET:HB2	1:A:810:PRO:HG2	1.77	0.66
1:A:272:LEU:HB3	1:A:305:VAL:HG11	1.76	0.66
1:A:662:GLN:OE1	1:A:1030:LEU:HD22	1.95	0.66
1:A:382:PHE:HE2	1:A:398:ARG:HD3	1.61	0.66
1:A:651:LEU:HD22	1:A:655:ASP:HB3	1.77	0.65
1:A:597:LYS:HB2	1:A:603:ILE:HD12	1.77	0.65
1:A:278:ASP:HB2	1:A:784:ARG:HH12	1.62	0.64
1:A:881:ILE:HG23	2:A:1201:O5E:H30	1.79	0.64
1:A:625:GLY:O	1:A:629:GLN:HG3	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:600:GLN:HG2	1:A:603:ILE:HG13	1.79	0.63
1:A:579:ARG:HD2	1:A:610:LEU:HD11	1.80	0.62
1:A:466:LEU:HD11	1:A:476:ARG:HD3	1.82	0.62
1:A:386:ASN:HB2	1:A:430:ASN:HB3	1.81	0.62
1:A:429:LEU:HB2	1:A:468:LEU:HD21	1.83	0.61
1:A:410:TRP:HB3	1:A:412:VAL:HG12	1.85	0.58
1:A:364:LYS:NZ	1:A:411:ASN:OD1	2.37	0.57
1:A:890:LYS:HG3	1:A:893:GLN:HE21	1.70	0.57
1:A:983:VAL:HG13	1:A:1082:VAL:HG21	1.88	0.56
1:A:1040:PRO:O	1:A:1042:LEU:HD13	2.04	0.56
1:A:1032:SER:HA	1:A:1048:ILE:HD12	1.87	0.56
1:A:374:PRO:C	1:A:376:ASN:HA	2.26	0.56
1:A:1084:PHE:O	1:A:1088:LEU:HD13	2.06	0.56
1:A:835:GLY:HA2	1:A:875:LYS:O	2.05	0.55
1:A:397:ARG:NH2	1:A:417:SER:OG	2.37	0.55
1:A:304:HIS:HB2	1:A:823:LEU:HD21	1.88	0.55
1:A:424:PRO:HD2	1:A:427:ALA:HB2	1.89	0.54
1:A:483:HIS:CD2	1:A:510:LYS:HG2	2.42	0.54
1:A:767:LEU:HD22	1:A:803:VAL:HG23	1.88	0.54
1:A:726:THR:HA	1:A:729:LEU:HD12	1.88	0.54
1:A:384:GLU:OE2	1:A:398:ARG:NE	2.41	0.54
1:A:929:VAL:HG22	1:A:995:MET:HG2	1.90	0.54
1:A:424:PRO:HG3	1:A:598:TRP:O	2.09	0.53
1:A:605:ALA:O	1:A:609:GLN:HG3	2.09	0.53
1:A:374:PRO:O	1:A:376:ASN:HA	2.10	0.52
1:A:939:THR:HB	1:A:945:GLY:HA2	1.92	0.52
1:A:509:ASP:OD2	1:A:512:ASN:HB2	2.10	0.51
1:A:911:LEU:O	1:A:915:SER:OG	2.23	0.51
1:A:890:LYS:HA	1:A:893:GLN:CD	2.31	0.51
1:A:905:GLU:HG2	1:A:993:PHE:CZ	2.46	0.51
1:A:983:VAL:HG22	1:A:1082:VAL:HG11	1.92	0.51
1:A:855:TRP:HZ3	1:A:864:LEU:HD11	1.76	0.50
1:A:475:LEU:HD21	1:A:522:ASN:HB2	1.93	0.50
1:A:357:CYS:SG	1:A:359:ARG:HG3	2.52	0.50
1:A:735:GLN:O	1:A:739:ILE:HG12	2.11	0.50
1:A:217:ASN:OD1	1:A:217:ASN:O	2.30	0.49
1:A:806:SER:O	1:A:806:SER:OG	2.30	0.49
1:A:1043:THR:HG23	1:A:1047:ASP:H	1.78	0.49
1:A:777:SER:HB3	1:A:778:GLN:HB2	1.94	0.49
1:A:890:LYS:HA	1:A:893:GLN:NE2	2.28	0.49
1:A:1011:ASP:OD1	1:A:1015:LYS:HE3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:VAL:O	1:A:185:MET:HG3	2.12	0.48
1:A:576:TRP:CZ2	1:A:603:ILE:HG23	2.48	0.48
1:A:862:LEU:HB3	1:A:934:GLY:HA3	1.94	0.48
1:A:893:GLN:O	1:A:896:VAL:O	2.32	0.48
1:A:890:LYS:HA	1:A:893:GLN:HG2	1.96	0.47
1:A:498:ASN:OD1	1:A:499:ALA:N	2.47	0.47
1:A:552:ARG:HH22	1:A:581:GLU:CD	2.17	0.47
1:A:552:ARG:HH12	1:A:581:GLU:CG	2.27	0.47
1:A:880:GLU:O	2:A:1201:OSE:H28	2.14	0.47
1:A:477:ARG:HA	1:A:520:LEU:HB3	1.98	0.46
1:A:995:MET:O	1:A:1005:HIS:HB2	2.15	0.46
1:A:547:MET:HB2	1:A:552:ARG:HH21	1.81	0.46
1:A:645:VAL:HA	1:A:648:LEU:HD12	1.98	0.46
1:A:177:ARG:NH2	1:A:718:GLU:OE2	2.49	0.46
1:A:706:SER:O	1:A:710:GLN:HB3	2.16	0.46
1:A:526:PRO:HB2	1:A:527:ILE:H	1.50	0.45
1:A:378:ASP:OD1	1:A:378:ASP:N	2.50	0.45
1:A:221:PHE:O	1:A:222:ILE:HD13	2.17	0.45
1:A:278:ASP:OD2	1:A:784:ARG:NH2	2.50	0.45
1:A:731:ASP:O	1:A:735:GLN:HG3	2.16	0.45
1:A:775:GLN:OE1	1:A:798:ILE:HD11	2.15	0.45
1:A:565:ASN:HA	1:A:566:PRO:HD3	1.77	0.45
1:A:899:THR:HB	1:A:901:ALA:H	1.82	0.44
1:A:282:VAL:HG22	1:A:283:GLY:N	2.31	0.44
1:A:792:LYS:HE3	1:A:792:LYS:HB3	1.66	0.44
1:A:651:LEU:HD22	1:A:655:ASP:CB	2.47	0.44
1:A:735:GLN:NE2	1:A:784:ARG:HB2	2.33	0.44
1:A:561:THR:O	1:A:591:LYS:NZ	2.45	0.44
1:A:853:SER:O	1:A:856:GLU:HB3	2.18	0.44
1:A:614:ARG:HG2	1:A:617:TRP:HB3	2.00	0.44
1:A:166:SER:O	1:A:510:LYS:NZ	2.47	0.43
1:A:774:LEU:HB3	1:A:779:LEU:HD23	1.99	0.43
1:A:990:ASP:O	1:A:994:VAL:HG23	2.17	0.43
1:A:625:GLY:HA2	1:A:1026:LEU:HD23	2.00	0.43
1:A:500:ASP:HB3	1:A:708:HIS:CD2	2.53	0.43
1:A:1032:SER:HA	1:A:1048:ILE:CD1	2.49	0.43
1:A:774:LEU:HD23	1:A:774:LEU:HA	1.85	0.43
1:A:865:LEU:HD21	1:A:882:VAL:HG11	1.99	0.43
1:A:787:TYR:CE1	1:A:880:GLU:HB2	2.54	0.43
1:A:368:ILE:HG23	1:A:410:TRP:HZ3	1.84	0.43
1:A:600:GLN:HB2	1:A:600:GLN:HE21	1.74	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:376:ASN:HB3	1:A:377:THR:C	2.39	0.42
1:A:1035:LEU:HD12	1:A:1048:ILE:CD1	2.35	0.42
1:A:278:ASP:CB	1:A:784:ARG:HH12	2.31	0.42
1:A:963:ILE:HD12	2:A:1201:OSE:C19	2.50	0.42
1:A:198:MET:SD	1:A:282:VAL:HG11	2.60	0.42
1:A:697:TRP:HZ2	1:A:735:GLN:HB3	1.84	0.41
1:A:182:THR:HB	1:A:183:PRO:HD3	2.02	0.41
1:A:498:ASN:OD1	1:A:1036:MET:O	2.37	0.41
1:A:768:LYS:O	1:A:772:GLU:HG3	2.20	0.41
1:A:389:HIS:O	1:A:392:GLN:HB3	2.21	0.41
1:A:739:ILE:O	1:A:743:GLN:HG3	2.21	0.41
1:A:607:THR:O	1:A:610:LEU:HB2	2.20	0.41
1:A:476:ARG:HG2	1:A:520:LEU:HD22	2.02	0.41
1:A:557:ALA:O	1:A:561:THR:HG23	2.20	0.41
1:A:857:THR:OG1	1:A:858:GLU:OE1	2.36	0.41
1:A:147:SER:HA	1:A:319:ARG:NH2	2.36	0.41
1:A:1058:GLY:O	1:A:1059:LYS:HD2	2.21	0.41
1:A:702:GLU:O	1:A:706:SER:HB3	2.21	0.40
1:A:308:ASP:OD1	1:A:308:ASP:N	2.54	0.40
1:A:997:THR:OG1	1:A:1001:LYS:O	2.28	0.40
1:A:1084:PHE:CZ	1:A:1088:LEU:HD11	2.57	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	816/966 (84%)	764 (94%)	45 (6%)	7 (1%)	21 56

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	807	LYS
1	A	1040	PRO
1	A	526	PRO
1	A	1045	LYS
1	A	862	LEU
1	A	897	GLY
1	A	371	PRO

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	744/864 (86%)	724 (97%)	20 (3%)	52 83

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	184	ARG
1	A	229	THR
1	A	477	ARG
1	A	600	GLN
1	A	638	GLU
1	A	744	LYS
1	A	777	SER
1	A	791	LEU
1	A	806	SER
1	A	823	LEU
1	A	832	PHE
1	A	907	LEU
1	A	959	ASN
1	A	983	VAL
1	A	1026	LEU
1	A	1032	SER
1	A	1039	MET
1	A	1076	ARG
1	A	1088	LEU
1	A	1089	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	483	HIS
1	A	600	GLN
1	A	825	ASN
1	A	959	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](#)

1 ligand is 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	0SE	A	1201	-	37,38,38	0.89	2 (5%)	54,55,55	2.18	20 (37%)

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	0SE	A	1201	-	-	0/18/36/36	0/5/5/5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1201	0SE	C14-N20	-2.03	1.34	1.37
2	A	1201	0SE	C24-N27	2.32	1.38	1.34

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1201	0SE	N20-C19-N18	-4.81	123.10	126.20
2	A	1201	0SE	C5-C6-N7	-3.98	104.45	110.96
2	A	1201	0SE	C21-C26-N25	-3.37	118.89	124.34
2	A	1201	0SE	C21-C22-N23	-3.25	119.08	124.34
2	A	1201	0SE	C6-C5-C4	-3.03	103.98	110.79
2	A	1201	0SE	C15-C17-N18	-2.97	117.41	120.48
2	A	1201	0SE	N25-C24-N23	-2.66	121.90	124.58
2	A	1201	0SE	C12-C13-C14	-2.58	118.08	120.88
2	A	1201	0SE	C5-C4-C2	-2.51	108.78	114.46
2	A	1201	0SE	C12-C11-N16	-2.45	120.27	123.01
2	A	1201	0SE	C29-N28-C17	-2.05	112.45	117.56
2	A	1201	0SE	N27-C24-N25	2.02	119.22	117.39
2	A	1201	0SE	C11-N16-C15	2.19	122.25	117.62
2	A	1201	0SE	C13-C14-N20	2.20	122.22	118.73
2	A	1201	0SE	C17-N18-C19	2.83	121.52	116.38
2	A	1201	0SE	N18-C17-N28	2.85	121.34	116.79
2	A	1201	0SE	C26-C21-C22	3.02	118.40	114.25
2	A	1201	0SE	C22-N23-C24	3.49	120.86	116.05
2	A	1201	0SE	C26-N25-C24	3.49	120.87	116.05
2	A	1201	0SE	C19-N20-C14	7.23	121.13	116.21

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1201	0SE	3	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	834/966 (86%)	0.29	56 (6%) 21 16	42, 96, 169, 235	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1041	GLN	9.6
1	A	375	ARG	8.4
1	A	376	ASN	7.2
1	A	1090	LEU	6.2
1	A	898	ASN	6.1
1	A	374	PRO	5.2
1	A	1091	VAL	5.1
1	A	1000	LYS	4.3
1	A	1084	PHE	4.2
1	A	932	CYS	3.8
1	A	1042	LEU	3.7
1	A	1046	GLU	3.6
1	A	226	ARG	3.6
1	A	527	ILE	3.5
1	A	823	LEU	3.5
1	A	1044	SER	3.5
1	A	526	PRO	3.4
1	A	211	LEU	3.4
1	A	230	SER	3.2
1	A	986	VAL	3.2
1	A	373	LEU	3.0
1	A	215	ILE	3.0
1	A	143	MET	3.0
1	A	1092	LEU	2.9
1	A	220	ILE	2.9
1	A	1087	PHE	2.9
1	A	489	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	895	THR	2.8
1	A	522	ASN	2.7
1	A	916	PRO	2.7
1	A	378	ASP	2.7
1	A	267	GLU	2.7
1	A	1045	LYS	2.7
1	A	636	SER	2.7
1	A	221	PHE	2.7
1	A	901	ALA	2.7
1	A	553	LYS	2.7
1	A	897	GLY	2.6
1	A	217	ASN	2.5
1	A	377	THR	2.4
1	A	776	ASN	2.4
1	A	216	ALA	2.3
1	A	936	CYS	2.3
1	A	909	HIS	2.3
1	A	320	LYS	2.3
1	A	919	GLU	2.2
1	A	1088	LEU	2.2
1	A	231	GLN	2.1
1	A	730	HIS	2.1
1	A	824	SER	2.1
1	A	148	GLN	2.1
1	A	270	PHE	2.1
1	A	269	ASP	2.1
1	A	899	THR	2.0
1	A	646	GLN	2.0
1	A	417	SER	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	0SE	A	1201	34/34	0.88	0.25	1.15	80,91,120,124	0

6.5 Other polymers [i](#)

There are no such residues in this entry.