



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:20 AM GMT

PDB ID : 3I25
Title : Potent Beta-Secretase 1 hydroxyethylene Inhibitor
Authors : Lindberg, J.D.; Borkakoti, N.; Nystrom, S.
Deposited on : 2009-06-29
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

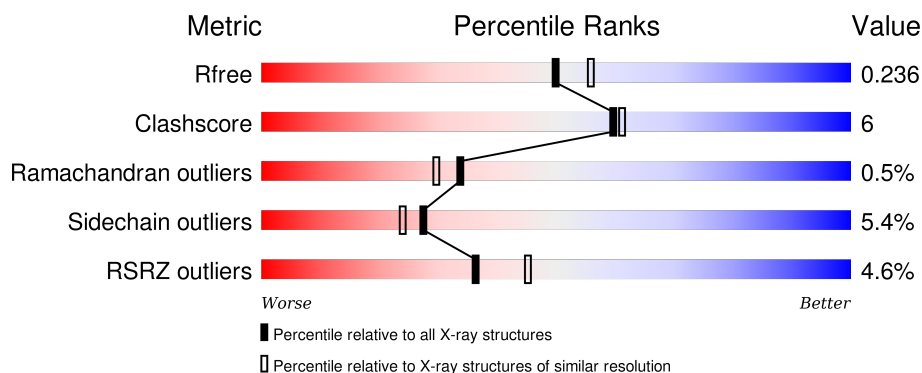
1 Overall quality at a glance

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	406	<div> <div>5%</div> <div> <div></div> <div>77%</div> <div>13%</div> <div>• 8%</div> </div> </div>
1	B	406	<div> <div>3%</div> <div> <div></div> <div>78%</div> <div>13%</div> <div>• 8%</div> </div> </div>
1	C	406	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>12%</div> <div>• 8%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9203 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.

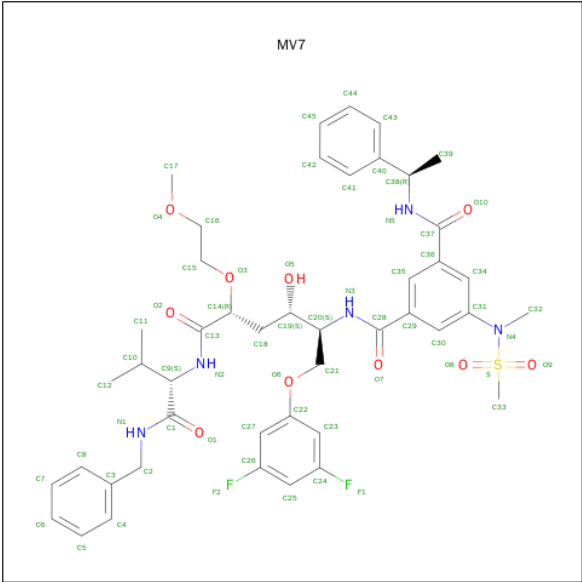
- Molecule 1 is a protein called Beta-secretase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	373	Total	C	N	O	S	0	0	0
			2940	1882	489	555	14			
1	B	373	Total	C	N	O	S	0	0	0
			2940	1882	489	555	14			
1	C	373	Total	C	N	O	S	0	0	0
			2940	1882	489	555	14			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	28P	MET	-	INITIATING METHIONINE	UNP P56817
B	28P	MET	-	INITIATING METHIONINE	UNP P56817
C	28P	MET	-	INITIATING METHIONINE	UNP P56817

- Molecule 2 is N-[(2S,3S,5R)-1-(3,5-DIFLUOROPHENOXY)-3-HYDROXY-5-(2-METHOXYETHOXY)-6-[[[(2S)-3-METHYL-1-OXO-1-(PHENYLMETHYLAMINO)BUTAN-2-YL]AMINO]-6-OXO-HEXAN-2-YL]-5-(METHYL-METHYLSULFONYL-AMINO)-N'-[(1R)-1-PHENYLETHYL]BENZENE-1,3-DICARBOXAMIDE (three-letter code: MV7) (formula: C₄₅H₅₅F₂N₅O₁₀S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	S	0	0
			63	45	2	5	10	1		
2	B	1	Total	C	F	N	O	S	0	0
			63	45	2	5	10	1		
2	C	1	Total	C	F	N	O	S	0	0
			63	45	2	5	10	1		

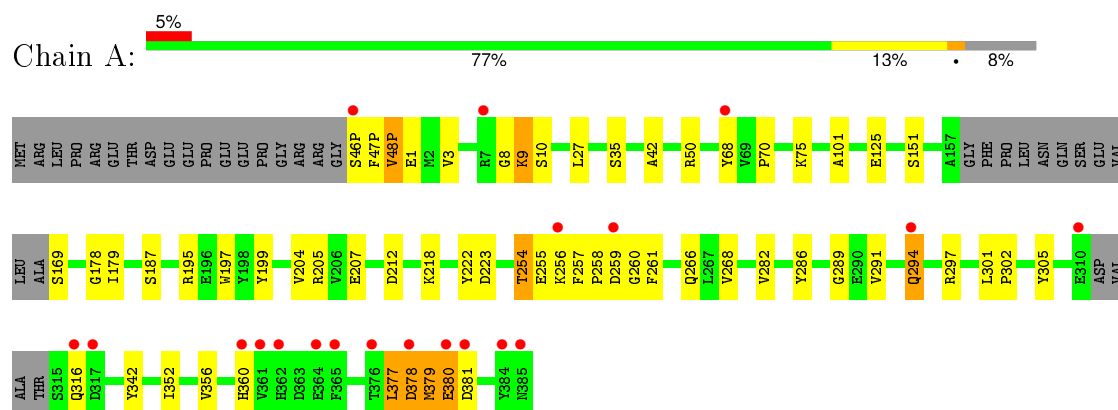
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	69	Total	O	0	0
			69	69		
3	B	67	Total	O	0	0
			67	67		
3	C	58	Total	O	0	0
			58	58		

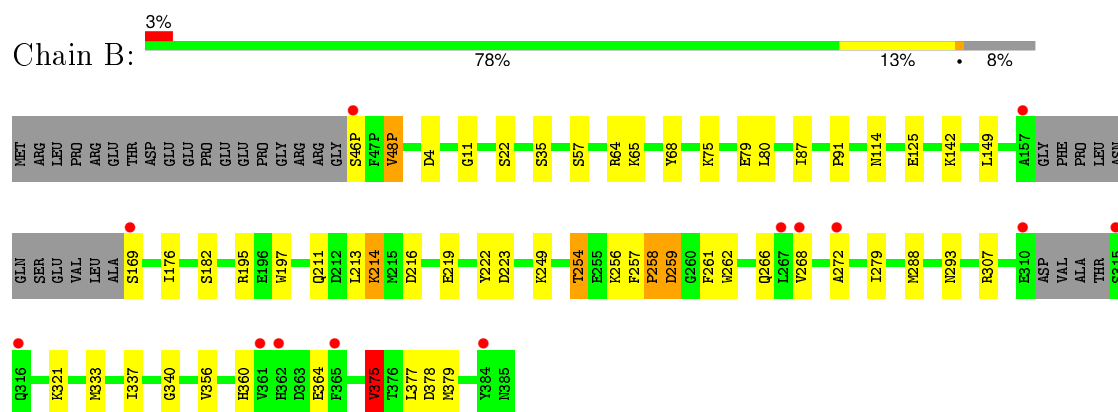
3 Residue-property plots [i](#)

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.

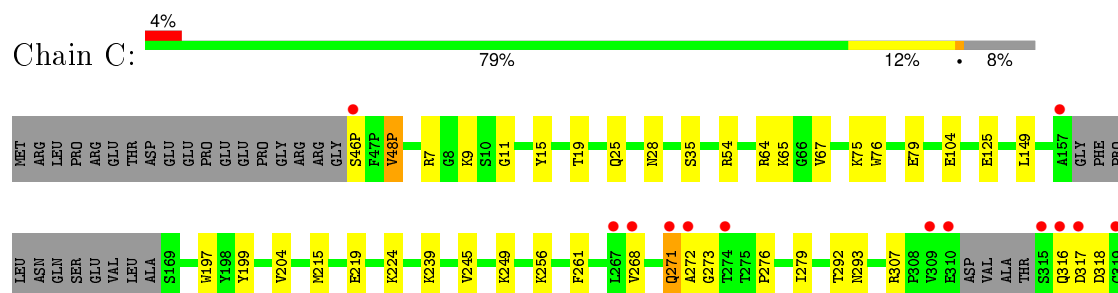
• Molecule 1: Beta-secretase 1

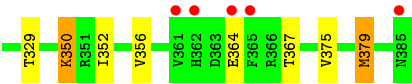


• Molecule 1: Beta-secretase 1



• Molecule 1: Beta-secretase 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.42Å 102.48Å 100.19Å 90.00° 103.53° 90.00°	Depositor
Resolution (Å)	79.16 – 2.10 79.16 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.7 (79.16-2.10) 95.7 (79.16-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.200 , 0.238 0.200 , 0.236	Depositor DCC
R_{free} test set	4449 reflections (5.24%)	DCC
Wilson B-factor (Å ²)	36.7	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 39.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 89314 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9203	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MV7

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	0.89	0/3014	0.85	2/4092 (0.0%)
1	B	0.91	0/3014	0.84	4/4092 (0.1%)
1	C	0.86	0/3014	0.82	1/4092 (0.0%)
All	All	0.89	0/9042	0.83	7/12276 (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	B	0	1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	48(P)	VAL	CB-CA-C	-6.51	99.04	111.40
1	B	375	VAL	CB-CA-C	-6.37	99.29	111.40
1	B	48(P)	VAL	CB-CA-C	-5.77	100.43	111.40
1	A	48(P)	VAL	CB-CA-C	-5.48	100.99	111.40
1	B	149	LEU	CA-CB-CG	5.38	127.69	115.30
1	A	212	ASP	CB-CG-OD1	5.30	123.07	118.30
1	B	375	VAL	CG1-CB-CG2	5.12	119.10	110.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	340	GLY	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	2940	0	2849	44	0
1	B	2940	0	2849	29	1
1	C	2940	0	2849	28	1
2	A	63	0	55	1	0
2	B	63	0	55	2	0
2	C	63	0	55	6	0
3	A	69	0	0	0	0
3	B	67	0	0	1	0
3	C	58	0	0	1	0
All	All	9203	0	8712	101	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 6.

All (101) 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:378:ASP:HB3	1:A:380:GLU:OE2	1.38	1.19
1:A:289:GLY:HA3	1:A:294:GLN:HG3	1.31	1.11
1:A:378:ASP:CB	1:A:380:GLU:OE2	2.08	1.01
1:A:289:GLY:HA3	1:A:294:GLN:CG	1.93	0.98
1:C:271:GLN:HA	1:C:317:ASP:HB2	1.50	0.93
1:C:204:VAL:HG11	1:C:379:MET:HG2	1.52	0.91
1:C:224:LYS:HE2	2:C:531:MV7:H17	1.55	0.86
1:C:215:MET:HE1	1:C:239:LYS:HG2	1.64	0.79
1:A:378:ASP:C	1:A:380:GLU:H	1.86	0.78
1:B:125:GLU:OE1	1:B:195:ARG:NH2	2.20	0.74
1:C:271:GLN:O	1:C:273:GLY:N	2.21	0.73
1:A:289:GLY:CA	1:A:294:GLN:HG3	2.17	0.73
1:A:378:ASP:C	1:A:380:GLU:N	2.38	0.73
1:A:257:PHE:HD2	1:A:268:VAL:HG11	1.55	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:257:PHE:HD2	1:B:268:VAL:HG11	1.55	0.70
1:A:8:GLY:C	1:A:9:LYS:HE2	2.16	0.66
1:A:378:ASP:CA	1:A:380:GLU:OE2	2.45	0.65
1:A:378:ASP:HA	1:A:380:GLU:OE2	1.98	0.64
1:A:291:VAL:O	1:A:294:GLN:HB3	1.98	0.63
1:A:260:GLY:HA3	1:A:266:GLN:HE21	1.63	0.63
1:A:9:LYS:N	1:A:9:LYS:HE2	2.13	0.62
1:B:211:GLN:HG3	3:B:400:HOH:O	1.99	0.62
1:A:378:ASP:O	1:A:380:GLU:N	2.33	0.61
1:A:257:PHE:CD2	1:A:268:VAL:HG11	2.36	0.60
1:B:261:PHE:CE1	1:B:268:VAL:HG22	2.37	0.60
1:C:9:LYS:HA	1:C:9:LYS:HE2	1.84	0.60
1:B:258:PRO:HB3	1:B:266:GLN:NE2	2.18	0.59
1:C:224:LYS:CE	2:C:531:MV7:H17	2.31	0.58
1:A:199:TYR:HB3	1:A:352:ILE:HD11	1.86	0.58
1:B:257:PHE:CD2	1:B:268:VAL:HG11	2.38	0.57
1:A:205:ARG:NH1	1:A:207:GLU:OE2	2.38	0.56
1:A:125:GLU:OE2	1:A:195:ARG:NH2	2.38	0.56
1:A:204:VAL:HG11	1:A:379:MET:HG2	1.89	0.54
1:A:291:VAL:HG22	1:A:294:GLN:HB2	1.89	0.54
1:A:377:LEU:HD13	1:A:377:LEU:N	2.23	0.53
1:C:215:MET:HE2	1:C:219:GLU:HG3	1.90	0.52
1:B:11:GLY:HA3	1:B:307:ARG:HH21	1.73	0.52
1:C:125:GLU:HG3	1:C:125:GLU:O	2.10	0.51
1:A:47(P):PHE:CZ	1:A:178:GLY:HA3	2.46	0.51
1:A:289:GLY:HA3	1:A:294:GLN:HG2	1.89	0.51
1:C:15:TYR:CD1	1:C:28:ASN:HB3	2.45	0.51
1:B:11:GLY:HA3	1:B:307:ARG:NH2	2.27	0.50
1:B:65:LYS:HG2	1:B:80:LEU:HD12	1.94	0.50
1:B:11:GLY:CA	1:B:307:ARG:HH21	2.25	0.49
1:B:91:PRO:HD3	1:B:176:ILE:HB	1.94	0.49
1:B:258:PRO:O	1:B:259:ASP:HB3	2.13	0.49
1:A:254:THR:HG22	1:A:255:GLU:HG3	1.93	0.48
1:B:125:GLU:OE1	1:B:195:ARG:NE	2.46	0.48
1:B:259:ASP:OD2	1:B:266:GLN:NE2	2.46	0.48
1:C:292:THR:O	1:C:293:ASN:HB2	2.13	0.48
1:A:261:PHE:CD1	1:A:268:VAL:HG13	2.49	0.48
1:A:27:LEU:HD23	1:A:50:ARG:HG3	1.95	0.48
1:C:19:THR:HA	1:C:25:GLN:O	2.14	0.47
1:A:68:TYR:OH	1:A:70:PRO:HB3	2.14	0.47
1:B:35:SER:HA	2:B:531:MV7:H11B	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:380:GLU:HG3	1:A:380:GLU:H	1.28	0.46
1:A:9:LYS:O	1:A:10:SER:C	2.54	0.46
1:A:35:SER:HA	2:A:531:MV7:H11B	1.97	0.46
1:A:302:PRO:HA	1:A:305:TYR:CE2	2.51	0.45
1:C:219:GLU:OE1	1:C:239:LYS:HE3	2.16	0.45
1:B:213:LEU:O	1:B:214:LYS:CB	2.65	0.45
1:C:329:THR:CB	2:C:531:MV7:H17B	2.47	0.45
1:B:257:PHE:O	1:B:258:PRO:C	2.55	0.45
1:C:329:THR:HB	2:C:531:MV7:H17B	1.98	0.45
1:C:379:MET:HB2	3:C:399:HOH:O	2.16	0.44
1:B:216:ASP:HB3	1:B:219:GLU:HG2	1.98	0.44
1:C:276:PRO:HB2	1:C:279:ILE:HG12	2.00	0.44
1:B:288:MET:HE1	1:B:379:MET:HA	1.99	0.43
1:A:282:VAL:HG12	1:A:301:LEU:HD23	2.00	0.43
1:C:261:PHE:CD1	1:C:268:VAL:HG23	2.54	0.43
1:B:222:TYR:HA	1:B:223:ASP:HA	1.86	0.42
2:B:531:MV7:H17B	2:B:531:MV7:H15A	1.81	0.42
1:C:67:VAL:HG13	1:C:76:TRP:HZ3	1.85	0.42
1:C:11:GLY:HA2	1:C:307:ARG:HH12	1.84	0.42
1:A:378:ASP:HA	1:A:380:GLU:HG3	2.01	0.42
1:B:288:MET:HE2	1:B:378:ASP:HA	2.02	0.42
1:C:199:TYR:HB3	1:C:352:ILE:HD11	2.02	0.42
1:C:329:THR:OG1	2:C:531:MV7:H17B	2.20	0.42
1:B:321:LYS:HB2	1:B:321:LYS:HE3	1.88	0.42
1:A:294:GLN:HE21	1:A:294:GLN:HB2	1.62	0.42
1:B:333:MET:HE3	1:B:337:ILE:HG21	2.02	0.42
1:B:258:PRO:O	1:B:259:ASP:CB	2.68	0.41
1:A:218:LYS:HG3	1:A:381:ASP:O	2.19	0.41
1:B:293:ASN:HA	1:B:375:VAL:HA	2.02	0.41
1:A:222:TYR:HA	1:A:223:ASP:HA	1.73	0.41
1:B:254:THR:HG21	1:B:279:ILE:HG21	2.02	0.41
1:C:350:LYS:HB2	1:C:350:LYS:HE2	1.90	0.41
1:A:261:PHE:CE1	1:A:268:VAL:HG22	2.55	0.41
1:C:9:LYS:CA	1:C:9:LYS:HE2	2.48	0.41
1:A:3:VAL:HG11	1:B:68:TYR:CZ	2.55	0.41
1:A:257:PHE:HA	1:A:258:PRO:HD2	1.91	0.41
1:C:54:ARG:HH12	1:C:79:GLU:CD	2.24	0.41
1:A:286:TYR:CZ	1:A:297:ARG:HD3	2.56	0.41
1:C:245:VAL:HG12	1:C:249:LYS:HD3	2.04	0.40
1:C:7:ARG:HD3	1:C:7:ARG:HA	1.91	0.40
1:C:35:SER:HA	2:C:531:MV7:H11B	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:179:ILE:HG23	1:A:342:TYR:HE2	1.86	0.40
1:B:249:LYS:HG3	1:B:262:TRP:CZ2	2.56	0.40
1:B:22:SER:O	1:B:57:SER:HA	2.21	0.40
1:A:46(P):SER:O	1:A:1:GLU:HB2	2.21	0.40
1:A:42:ALA:CB	1:A:101:ALA:HB1	2.51	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)
1:B:46(P):SER:OG	1:C:104:GLU:OE2[2_455]	2.13	0.07

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	367/406 (90%)	358 (98%)	8 (2%)	1 (0%)	46	45
1	B	367/406 (90%)	355 (97%)	9 (2%)	3 (1%)	24	17
1	C	367/406 (90%)	357 (97%)	8 (2%)	2 (0%)	34	30
All	All	1101/1218 (90%)	1070 (97%)	25 (2%)	6 (0%)	34	30

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	258	PRO
1	C	272	ALA
1	B	259	ASP
1	C	318	ASP
1	A	379	MET
1	B	272	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	319/347 (92%)	302 (95%)	17 (5%)	28	25
1	B	319/347 (92%)	300 (94%)	19 (6%)	24	20
1	C	319/347 (92%)	303 (95%)	16 (5%)	30	27
All	All	957/1041 (92%)	905 (95%)	52 (5%)	27	24

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48(P)	VAL
1	A	9	LYS
1	A	75	LYS
1	A	151	SER
1	A	169	SER
1	A	187	SER
1	A	197	TRP
1	A	254	THR
1	A	256	LYS
1	A	259	ASP
1	A	294	GLN
1	A	316	GLN
1	A	356	VAL
1	A	360	HIS
1	A	377	LEU
1	A	378	ASP
1	A	380	GLU
1	B	48(P)	VAL
1	B	4	ASP
1	B	64	ARG
1	B	75	LYS
1	B	79	GLU
1	B	87	ILE
1	B	114	ASN
1	B	142	LYS
1	B	169	SER

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Mol	Chain	Res	Type
1	B	182	SER
1	B	197	TRP
1	B	214	LYS
1	B	254	THR
1	B	256	LYS
1	B	356	VAL
1	B	360	HIS
1	B	364	GLU
1	B	375	VAL
1	B	377	LEU
1	C	46(P)	SER
1	C	48(P)	VAL
1	C	64	ARG
1	C	65	LYS
1	C	75	LYS
1	C	149	LEU
1	C	197	TRP
1	C	256	LYS
1	C	271	GLN
1	C	316	GLN
1	C	350	LYS
1	C	356	VAL
1	C	364	GLU
1	C	367	THR
1	C	375	VAL
1	C	379	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	266	GLN
1	A	294	GLN
1	A	304	GLN
1	B	28	ASN
1	B	266	GLN
1	B	304	GLN
1	B	316	GLN
1	B	326	GLN
1	B	360	HIS
1	C	28	ASN
1	C	293	ASN

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Mol	Chain	Res	Type
1	C	316	GLN
1	C	360	HIS
1	C	362	HIS

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	MV7	A	531	-	66,66,66	1.08	3 (4%)	81,91,91	1.80	15 (18%)
2	MV7	B	531	-	66,66,66	1.10	4 (6%)	81,91,91	1.59	14 (17%)
2	MV7	C	531	-	66,66,66	1.08	4 (6%)	81,91,91	1.66	18 (22%)

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	MV7	A	531	-	-	0/69/69/69	0/4/4/4
2	MV7	B	531	-	-	0/69/69/69	0/4/4/4
2	MV7	C	531	-	-	0/69/69/69	0/4/4/4

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	531	MV7	C33-S	-4.15	1.67	1.75
2	C	531	MV7	S-N4	-4.13	1.55	1.65
2	B	531	MV7	C33-S	-3.69	1.68	1.75
2	A	531	MV7	O9-S	-2.94	1.40	1.43
2	A	531	MV7	S-N4	-2.77	1.59	1.65
2	C	531	MV7	C33-S	-2.72	1.70	1.75
2	B	531	MV7	C31-N4	-2.18	1.41	1.44
2	C	531	MV7	C35-C36	-2.01	1.36	1.39
2	C	531	MV7	C27-C26	2.07	1.41	1.37
2	B	531	MV7	C23-C24	2.16	1.41	1.37
2	B	531	MV7	C25-C26	2.17	1.41	1.37

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	531	MV7	C10-C9-C1	-6.54	94.15	111.31
2	C	531	MV7	C10-C9-C1	-5.34	97.30	111.31
2	A	531	MV7	O9-S-C33	-5.27	101.41	108.70
2	A	531	MV7	C10-C9-C1	-5.16	97.77	111.31
2	A	531	MV7	C21-C20-N3	-4.97	103.02	109.60
2	C	531	MV7	O8-S-O9	-3.42	113.90	118.66
2	A	531	MV7	O8-S-O9	-3.38	113.95	118.66
2	C	531	MV7	O10-C37-C36	-2.67	116.42	120.97
2	C	531	MV7	O3-C15-C16	-2.62	100.25	109.68
2	B	531	MV7	O9-S-C33	-2.51	105.23	108.70
2	C	531	MV7	C3-C2-N1	-2.50	107.18	112.88
2	B	531	MV7	O2-C13-C14	-2.44	115.43	119.52
2	C	531	MV7	O8-S-C33	-2.41	105.36	108.70
2	B	531	MV7	O3-C14-C13	-2.36	106.22	111.47
2	A	531	MV7	C3-C2-N1	-2.21	107.83	112.88
2	A	531	MV7	C29-C35-C36	-2.21	118.05	120.47
2	C	531	MV7	C40-C38-N5	-2.12	107.48	111.41
2	B	531	MV7	C23-C24-C25	-2.07	120.78	123.52
2	C	531	MV7	C23-C24-C25	-2.07	120.79	123.52
2	A	531	MV7	C23-C24-C25	-2.02	120.85	123.52
2	C	531	MV7	C29-C35-C36	-2.01	118.27	120.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	531	MV7	O2-C13-N2	2.01	126.86	122.93
2	C	531	MV7	C10-C9-N2	2.06	116.78	111.36
2	C	531	MV7	C22-C23-C24	2.07	120.59	117.66
2	A	531	MV7	C35-C36-C34	2.11	122.15	119.64
2	C	531	MV7	O6-C21-C20	2.12	112.40	108.44
2	C	531	MV7	C2-N1-C1	2.17	125.33	122.34
2	B	531	MV7	C19-C20-N3	2.46	113.69	109.73
2	B	531	MV7	C11-C10-C12	2.48	117.96	110.67
2	A	531	MV7	O9-S-N4	2.51	110.19	106.94
2	A	531	MV7	O8-S-C33	2.58	112.26	108.70
2	B	531	MV7	O5-C19-C20	2.66	115.35	109.21
2	A	531	MV7	C2-N1-C1	2.66	126.00	122.34
2	C	531	MV7	C11-C10-C9	2.68	119.15	111.14
2	B	531	MV7	C10-C9-N2	2.95	119.15	111.36
2	C	531	MV7	C19-C20-N3	2.96	114.49	109.73
2	A	531	MV7	C11-C10-C9	3.00	120.12	111.14
2	B	531	MV7	O8-S-N4	3.00	110.84	106.94
2	B	531	MV7	C21-O6-C22	3.02	124.38	117.89
2	C	531	MV7	C35-C36-C34	3.04	123.26	119.64
2	A	531	MV7	O8-S-N4	3.15	111.02	106.94
2	B	531	MV7	C11-C10-C9	3.24	120.83	111.14
2	C	531	MV7	C21-O6-C22	3.28	124.95	117.89
2	A	531	MV7	C21-O6-C22	4.62	127.83	117.89
2	B	531	MV7	O6-C21-C20	5.03	117.85	108.44
2	A	531	MV7	O6-C21-C20	5.27	118.29	108.44
2	C	531	MV7	O8-S-N4	6.27	115.07	106.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	531	MV7	1	0
2	B	531	MV7	2	0
2	C	531	MV7	6	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, 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	373/406 (91%)	0.56	20 (5%) 29 38	19, 36, 59, 77	0
1	B	373/406 (91%)	0.48	13 (3%) 48 57	20, 35, 57, 73	0
1	C	373/406 (91%)	0.59	18 (4%) 34 43	19, 38, 70, 83	0
All	All	1119/1218 (91%)	0.54	51 (4%) 36 45	19, 36, 64, 83	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	315	SER	9.2
1	C	157	ALA	7.1
1	A	365	PHE	5.8
1	A	316	GLN	5.4
1	C	272	ALA	5.1
1	C	317	ASP	4.7
1	B	365	PHE	4.4
1	A	68	TYR	4.4
1	C	309	VAL	3.8
1	A	381	ASP	3.7
1	A	380	GLU	3.7
1	C	362	HIS	3.6
1	C	310	GLU	3.5
1	A	259	ASP	3.4
1	A	385	ASN	3.4
1	C	267	LEU	3.4
1	A	46(P)	SER	3.3
1	C	361	VAL	3.2
1	A	378	ASP	3.1
1	A	364	GLU	3.0
1	B	362	HIS	3.0
1	A	310	GLU	3.0
1	C	365	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	294	GLN	2.9
1	B	316	GLN	2.8
1	A	362	HIS	2.7
1	C	46(P)	SER	2.7
1	B	46(P)	SER	2.7
1	C	316	GLN	2.7
1	B	157	ALA	2.6
1	A	361	VAL	2.6
1	C	319	CYS	2.5
1	C	364	GLU	2.5
1	A	256	LYS	2.4
1	B	272	ALA	2.3
1	C	271	GLN	2.3
1	A	360	HIS	2.3
1	B	384	TYR	2.3
1	A	7	ARG	2.3
1	B	315	SER	2.2
1	B	310	GLU	2.2
1	A	317	ASP	2.2
1	A	384	TYR	2.2
1	B	268	VAL	2.1
1	C	385	ASN	2.1
1	B	169	SER	2.1
1	B	361	VAL	2.1
1	A	376	THR	2.1
1	C	268	VAL	2.1
1	B	267	LEU	2.0
1	C	274	THR	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. 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(Å ²)	Q<0.9
2	MV7	C	531	63/63	0.95	0.15	0.79	25,32,41,45	0
2	MV7	A	531	63/63	0.96	0.15	0.15	26,32,49,53	0
2	MV7	B	531	63/63	0.97	0.14	-0.17	22,28,40,46	0

6.5 Other polymers

There are no such residues in this entry.