



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:42 PM GMT

PDB ID : 4TYB  
Title : An Ligand-observed Mass Spectrometry-based Approach Integrated into the  
Fragment Based Lead Discovery Pipeline  
Authors : Shui, W.; Yang, C.; Lin, J.; Chen, X.; Qin, S.; Chen, S.  
Deposited on : 2014-07-08  
Resolution : 2.93 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](mailto: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.

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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

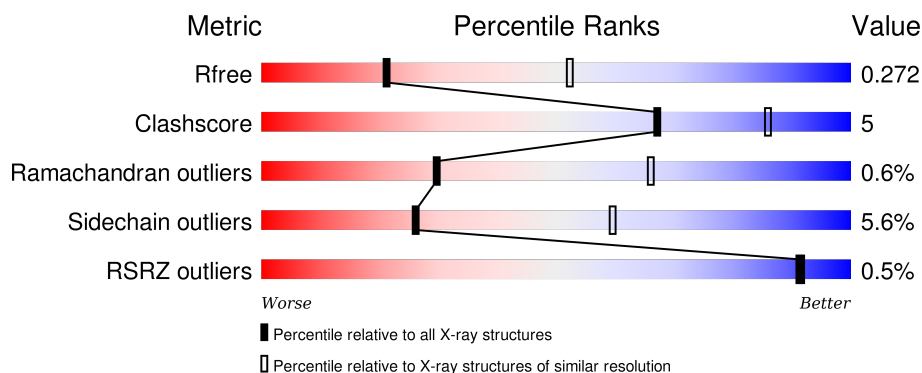
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*





The reported resolution of this entry is 2.93 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  $\geq 3$ , 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	566	 81% 13% . .
1	B	566	 82% 13% . .
1	C	566	 84% 11% . .
1	D	566	 82% 13% . .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	3B1	A	601	-	-	X	X
2	3B1	B	601	-	-	-	X
2	3B1	C	601	-	-	-	X
2	3B1	D	601	-	-	-	X

## 2 Entry composition [i](#)

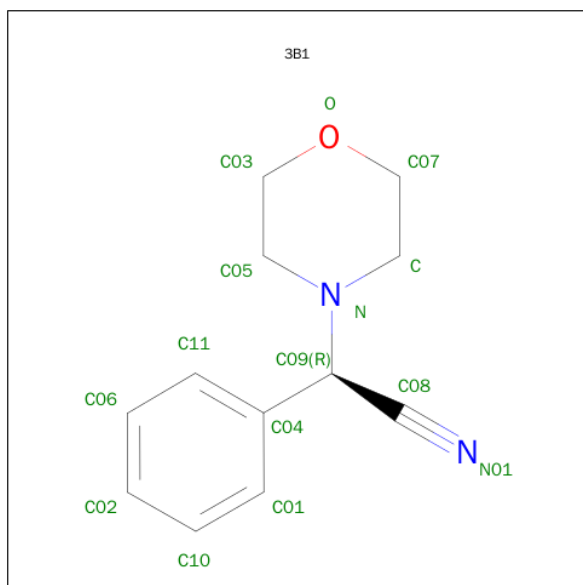
There are 2 unique types of molecules in this entry. The entry contains 16969 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Polypeptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	545	Total	C	N	O	S	0	0	0
			4229	2669	745	784	31			
1	B	544	Total	C	N	O	S	0	0	0
			4222	2664	744	783	31			
1	C	545	Total	C	N	O	S	0	0	0
			4229	2669	745	784	31			
1	D	545	Total	C	N	O	S	0	0	0
			4229	2669	745	784	31			

- Molecule 2 is (2R)-morpholin-4-yl(phenyl)ethanenitrile (three-letter code: 3B1) (formula:  $C_{12}H_{14}N_2O$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			15	12	2	1		
2	B	1	Total	C	N	O	0	0
			15	12	2	1		

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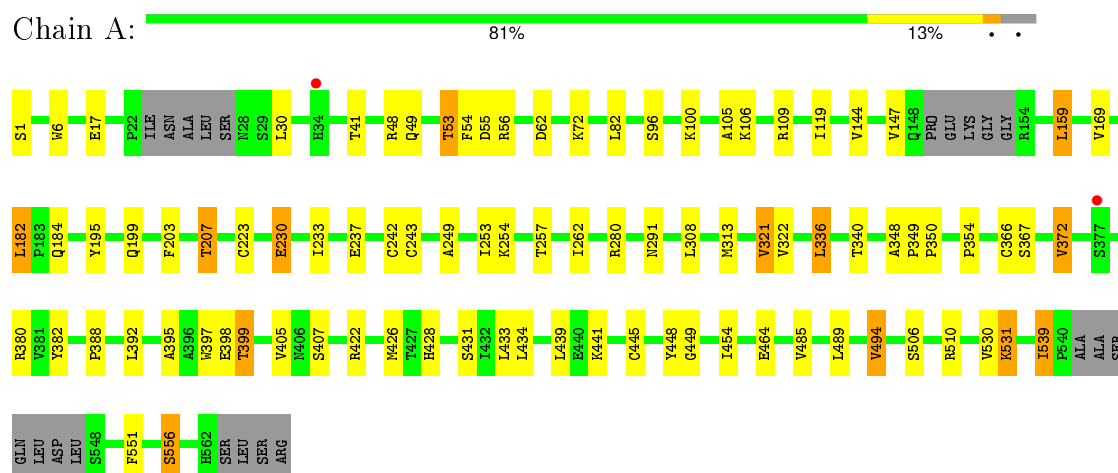
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			15	12	2	1		
2	D	1	Total	C	N	O	0	0
			15	12	2	1		

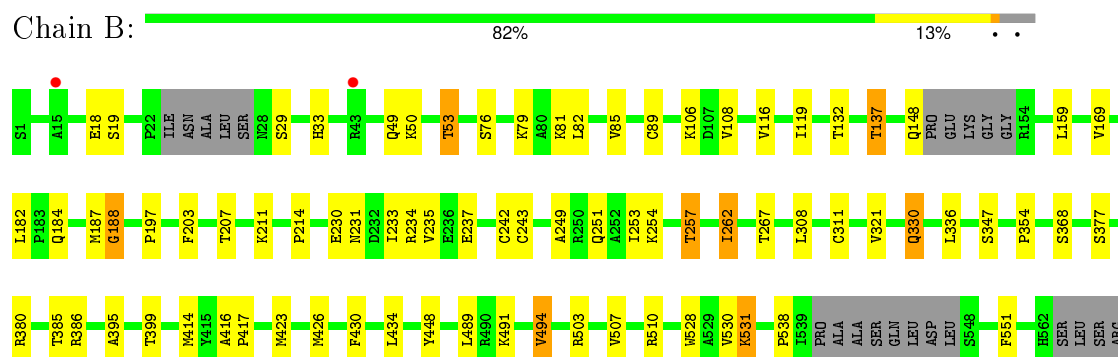
### 3 Residue-property plots

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

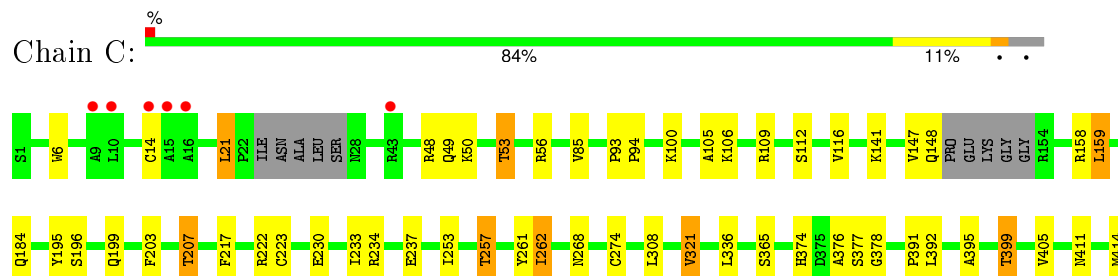
#### • Molecule 1: Polyprotein

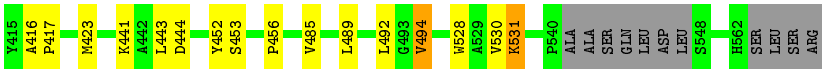


#### • Molecule 1: Polyprotein

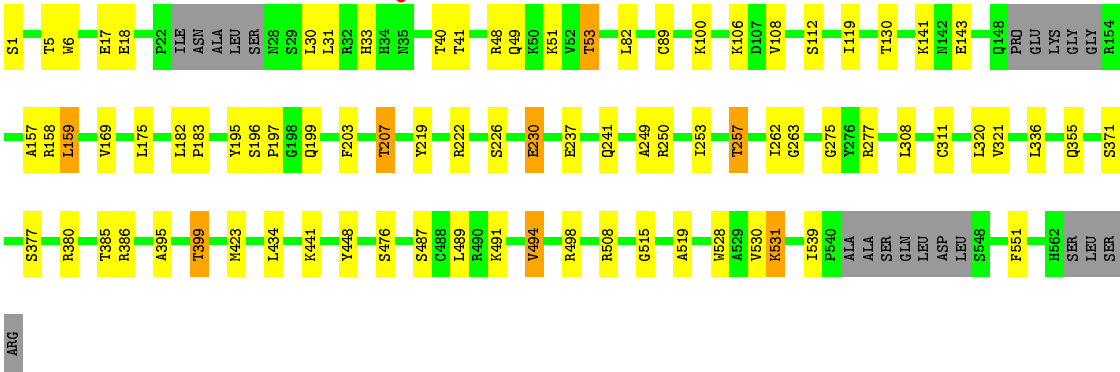
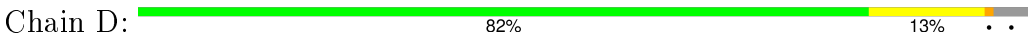


#### • Molecule 1: Polyprotein





● Molecule 1: Polyprotein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	102.45Å 102.35Å 251.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.41 – 2.93 45.78 – 2.93	Depositor EDS
% Data completeness (in resolution range)	98.6 (47.41-2.93) 98.6 (45.78-2.93)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.69 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.205 , 0.273 0.208 , 0.272	Depositor DCC
$R_{free}$ test set	2899 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.9	Xtriage
Anisotropy	0.065	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 14.7	EDS
Estimated twinning fraction	0.087 for k,h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 57097 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	16969	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.48% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 3B1

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	0.53	0/4321	0.73	1/5864 (0.0%)
1	B	0.53	0/4313	0.75	0/5852
1	C	0.53	0/4321	0.74	2/5864 (0.0%)
1	D	0.53	0/4321	0.74	1/5864 (0.0%)
All	All	0.53	0/17276	0.74	4/23444 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	321	VAL	CB-CA-C	-6.00	100.00	111.40
1	A	321	VAL	CB-CA-C	-5.47	101.00	111.40
1	D	321	VAL	CB-CA-C	-5.43	101.07	111.40
1	C	21	LEU	CA-CB-CG	5.38	127.68	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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	4229	0	4219	46	0
1	B	4222	0	4212	38	0
1	C	4229	0	4219	40	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	4229	0	4219	34	0
2	A	15	0	14	10	0
2	B	15	0	14	2	0
2	C	15	0	14	4	0
2	D	15	0	14	0	0
All	All	16969	0	16925	164	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 5.

All (164) 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:49:GLN:O	1:B:53:THR:HG22	1.55	1.06
1:D:49:GLN:O	1:D:53:THR:HG22	1.54	1.05
1:A:49:GLN:O	1:A:53:THR:HG23	1.62	0.99
1:A:366:CYS:HB3	2:A:601:3B1:H14	1.45	0.98
1:A:395:ALA:O	1:A:399:THR:HG22	1.72	0.89
1:D:203:PHE:O	1:D:207:THR:HG23	1.73	0.88
1:B:49:GLN:O	1:B:53:THR:CG2	2.26	0.83
1:C:253:ILE:O	1:C:257:THR:HG23	1.79	0.83
1:A:6:TRP:HE1	1:A:53:THR:HG21	1.43	0.81
1:A:366:CYS:HB3	2:A:601:3B1:C	2.11	0.79
1:C:203:PHE:O	1:C:207:THR:HG23	1.83	0.78
1:A:203:PHE:O	1:A:207:THR:HG23	1.83	0.78
1:B:395:ALA:O	1:B:399:THR:HG23	1.83	0.77
1:A:253:ILE:O	1:A:257:THR:HG23	1.85	0.77
1:C:49:GLN:O	1:C:53:THR:CG2	2.33	0.76
1:D:82:LEU:HD13	1:D:249:ALA:HB2	1.67	0.76
1:C:49:GLN:O	1:C:53:THR:HG23	1.89	0.72
1:D:49:GLN:O	1:D:53:THR:CG2	2.38	0.69
1:C:234:ARG:HG3	1:C:262:ILE:HD11	1.76	0.66
1:D:253:ILE:O	1:D:257:THR:HG23	1.97	0.64
1:A:366:CYS:SG	2:A:601:3B1:H11	2.38	0.64
1:A:340:THR:HG23	1:A:350:PRO:HG3	1.81	0.63
1:A:48:ARG:HG2	1:A:159:LEU:HD13	1.80	0.62
1:B:231:ASN:O	1:B:235:VAL:HG23	2.00	0.62
1:D:395:ALA:O	1:D:399:THR:HG22	2.02	0.60
2:A:601:3B1:H2	2:A:601:3B1:C	2.31	0.60
1:A:82:LEU:HD13	1:A:249:ALA:HB2	1.86	0.58
1:B:182:LEU:HD13	1:B:243:CYS:SG	2.43	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:601:3B1:H2	2:A:601:3B1:H15	1.85	0.57
1:A:366:CYS:CB	2:A:601:3B1:H14	2.26	0.57
1:B:233:ILE:HG22	1:B:262:ILE:HD12	1.87	0.57
1:D:31:LEU:O	1:D:31:LEU:HD12	2.06	0.56
1:A:449:GLY:HA3	1:A:556:SER:CA	2.36	0.55
1:D:308:LEU:HB3	1:D:311:CYS:SG	2.46	0.55
1:A:105:ALA:O	1:A:109:ARG:HG3	2.06	0.55
1:A:144:VAL:HG21	1:A:397:TRP:CG	2.42	0.55
1:B:182:LEU:CD1	1:B:243:CYS:SG	2.95	0.55
1:C:49:GLN:O	1:C:53:THR:HG22	2.07	0.54
1:A:336:LEU:HD12	1:A:354:PRO:HG2	1.87	0.54
1:A:237:GLU:HG3	1:A:257:THR:HG21	1.90	0.54
1:A:449:GLY:HA3	1:A:556:SER:HA	1.90	0.54
1:D:263:GLY:HA3	1:D:277:ARG:O	2.08	0.54
1:B:423:MET:HA	1:B:528:TRP:CZ2	2.43	0.54
1:B:253:ILE:O	1:B:257:THR:HG23	2.08	0.54
1:D:241:GLN:OE1	1:D:250:ARG:HG3	2.08	0.53
1:D:530:VAL:O	1:D:531:LYS:CB	2.57	0.53
2:A:601:3B1:H2	2:A:601:3B1:C07	2.39	0.53
1:B:503:ARG:O	1:B:507:VAL:HG23	2.09	0.53
1:A:489:LEU:HA	1:A:494:VAL:HG13	1.91	0.53
1:D:30:LEU:O	1:D:494:VAL:HB	2.09	0.52
1:D:237:GLU:HG3	1:D:257:THR:HG21	1.92	0.52
1:B:233:ILE:CG2	1:B:262:ILE:HD12	2.40	0.52
1:C:321:VAL:CG2	1:C:365:SER:CB	2.88	0.51
1:A:6:TRP:NE1	1:A:53:THR:HG21	2.21	0.51
1:B:85:VAL:HG11	1:B:116:VAL:HG13	1.91	0.51
1:B:434:LEU:HD21	1:B:510:ARG:HB2	1.91	0.51
1:B:203:PHE:O	1:B:207:THR:HG23	2.10	0.51
1:C:147:VAL:HG12	1:C:148:GLN:HG2	1.93	0.50
1:B:530:VAL:O	1:B:531:LYS:CB	2.59	0.50
1:A:464:GLU:OE2	1:A:539:ILE:HG12	2.10	0.50
1:B:426:MET:O	1:B:430:PHE:HB2	2.11	0.49
1:C:530:VAL:O	1:C:531:LYS:HB3	2.12	0.49
1:D:385:THR:OG1	1:D:386:ARG:N	2.46	0.49
1:B:237:GLU:CG	1:B:257:THR:HG21	2.42	0.49
1:A:280:ARG:NE	1:A:291:ASN:OD1	2.36	0.49
1:B:33:HIS:HD2	1:B:491:LYS:O	1.96	0.49
1:C:416:ALA:HB3	1:C:417:PRO:HD3	1.95	0.49
1:D:119:ILE:HD13	1:D:169:VAL:HG11	1.94	0.49
1:B:197:PRO:HG3	1:B:414:MET:CE	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:489:LEU:HD22	1:B:494:VAL:HG22	1.95	0.48
1:A:313:MET:HG2	1:A:322:VAL:HG22	1.95	0.48
1:D:17:GLU:OE1	1:D:41:THR:HB	2.13	0.48
1:A:434:LEU:HD21	1:A:510:ARG:CB	2.43	0.48
2:A:601:3B1:H10	2:A:601:3B1:H2	1.96	0.48
1:C:391:PRO:HB3	1:C:411:ASN:ND2	2.29	0.48
1:C:414:MET:HG3	2:C:601:3B1:C03	2.44	0.47
2:A:601:3B1:C01	2:A:601:3B1:H15	2.44	0.47
1:B:76:SER:HA	1:B:242:CYS:O	2.14	0.47
1:D:48:ARG:HG2	1:D:159:LEU:HD13	1.94	0.47
1:C:85:VAL:HG11	1:C:116:VAL:HG22	1.96	0.47
1:C:195:TYR:HA	1:C:199:GLN:HE21	1.79	0.47
1:C:6:TRP:HE1	1:C:53:THR:HG21	1.79	0.47
1:C:321:VAL:HG22	1:C:365:SER:HB3	1.98	0.46
1:C:423:MET:HA	1:C:528:TRP:CZ2	2.50	0.46
1:A:54:PHE:HD1	1:A:55:ASP:O	1.98	0.46
1:A:233:ILE:CG2	1:A:262:ILE:HD12	2.46	0.46
1:A:422:ARG:HA	1:A:426:MET:SD	2.55	0.46
1:D:515:GLY:HA2	1:D:519:ALA:HB2	1.97	0.46
1:B:330:GLN:HE21	1:B:330:GLN:HA	1.80	0.46
1:D:196:SER:H	1:D:199:GLN:HE21	1.63	0.46
1:D:40:THR:HB	1:D:157:ALA:HB2	1.96	0.46
1:B:336:LEU:HD12	1:B:354:PRO:HB2	1.98	0.46
1:C:233:ILE:CG2	1:C:262:ILE:HD12	2.46	0.45
1:B:197:PRO:HG3	1:B:414:MET:HE3	1.97	0.45
1:A:530:VAL:O	1:A:531:LYS:CB	2.64	0.45
1:A:372:VAL:HG12	1:A:382:TYR:CD1	2.50	0.45
1:A:434:LEU:HD21	1:A:510:ARG:HB2	1.98	0.45
1:B:187:MET:O	1:B:188:GLY:C	2.54	0.45
1:C:374:HIS:HD2	1:C:378:GLY:O	1.99	0.45
1:A:448:TYR:CE2	1:A:551:PHE:HD1	2.34	0.45
1:C:485:VAL:O	1:C:489:LEU:HG	2.16	0.45
1:C:444:ASP:HA	1:C:452:TYR:O	2.17	0.45
1:D:5:THR:O	1:D:275:GLY:HA3	2.17	0.44
1:A:428:HIS:O	1:A:431:SER:HB3	2.17	0.44
1:D:141:LYS:HE3	1:D:143:GLU:HG3	1.99	0.44
1:D:489:LEU:HA	1:D:494:VAL:HG13	1.99	0.44
1:A:348:ALA:N	1:A:349:PRO:CD	2.80	0.44
1:C:48:ARG:HG2	1:C:159:LEU:HD13	2.00	0.44
1:B:385:THR:OG1	1:B:386:ARG:N	2.50	0.44
1:B:251:GLN:HA	1:B:251:GLN:OE1	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:GLU:HB3	1:A:41:THR:HG22	2.00	0.44
1:D:6:TRP:HE1	1:D:53:THR:HG21	1.83	0.44
1:C:395:ALA:O	1:C:399:THR:HG23	2.17	0.43
1:B:448:TYR:CE2	1:B:551:PHE:HD1	2.36	0.43
1:C:257:THR:HA	1:C:261:TYR:HB2	1.99	0.43
1:D:448:TYR:CE2	1:D:551:PHE:HD1	2.36	0.43
1:C:443:LEU:HD12	1:C:456:PRO:HG3	1.99	0.43
1:B:230:GLU:O	1:B:234:ARG:HG3	2.19	0.43
1:A:388:PRO:O	1:A:392:LEU:HG	2.17	0.43
1:B:137:THR:HA	1:B:267:THR:O	2.19	0.43
1:C:237:GLU:CG	1:C:257:THR:HG21	2.49	0.43
1:D:423:MET:HA	1:D:528:TRP:CZ2	2.54	0.43
1:C:217:PHE:CD1	1:C:336:LEU:HD21	2.53	0.43
1:B:414:MET:HE2	2:B:601:3B1:H12	2.01	0.42
1:A:433:LEU:HB3	1:A:439:LEU:HD23	2.00	0.42
1:A:96:SER:HB3	1:A:105:ALA:HB2	2.01	0.42
1:C:414:MET:HG3	2:C:601:3B1:H5	2.01	0.42
1:B:416:ALA:HB3	1:B:417:PRO:HD3	2.01	0.42
1:D:219:TYR:HB3	1:D:320:LEU:HD23	2.01	0.42
1:D:33:HIS:HD2	1:D:491:LYS:O	2.02	0.42
1:C:489:LEU:HD22	1:C:494:VAL:CG2	2.49	0.42
1:A:119:ILE:HD13	1:A:169:VAL:HG11	2.02	0.42
1:C:321:VAL:CG2	1:C:365:SER:HB2	2.49	0.42
1:D:182:LEU:N	1:D:183:PRO:CD	2.82	0.42
1:A:72:LYS:HG2	1:A:242:CYS:SG	2.60	0.42
1:D:175:LEU:HG	1:D:253:ILE:HG12	2.01	0.42
1:D:263:GLY:CA	1:D:277:ARG:O	2.67	0.42
1:C:48:ARG:HG3	1:C:159:LEU:HD22	2.02	0.42
2:A:601:3B1:C01	2:A:601:3B1:C	2.96	0.42
1:B:530:VAL:O	1:B:531:LYS:HB3	2.20	0.42
1:D:89:CYS:SG	1:D:108:VAL:HG13	2.59	0.42
1:D:195:TYR:HA	1:D:199:GLN:HE21	1.85	0.42
1:C:268:ASN:HB3	1:C:274:CYS:SG	2.60	0.42
1:A:30:LEU:O	1:A:494:VAL:HB	2.20	0.41
1:C:321:VAL:HG22	1:C:365:SER:CB	2.49	0.41
1:A:182:LEU:HD13	1:A:243:CYS:SG	2.60	0.41
1:C:414:MET:CE	2:C:601:3B1:H6	2.51	0.41
1:A:1:SER:OG	1:A:230:GLU:OE1	2.30	0.41
1:C:414:MET:HE2	2:C:601:3B1:H12	2.01	0.41
1:B:119:ILE:HD13	1:B:169:VAL:HG11	2.02	0.41
1:A:485:VAL:O	1:A:489:LEU:HG	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:ALA:O	1:C:109:ARG:HG3	2.20	0.41
1:C:237:GLU:HG3	1:C:257:THR:HG21	2.01	0.41
1:D:196:SER:O	1:D:197:PRO:C	2.56	0.41
1:A:195:TYR:HA	1:A:199:GLN:HE21	1.86	0.41
1:C:321:VAL:CG2	1:C:365:SER:HB3	2.51	0.41
1:C:392:LEU:HD23	1:C:392:LEU:HA	1.94	0.41
1:B:211:LYS:HB2	1:B:214:PRO:HB3	2.02	0.40
1:A:398:GLU:OE2	1:A:407:SER:OG	2.28	0.40
1:A:445:CYS:SG	1:A:454:ILE:HD12	2.61	0.40
1:B:414:MET:HE2	2:B:601:3B1:H15	2.03	0.40
1:B:89:CYS:SG	1:B:108:VAL:HG13	2.62	0.40
1:C:93:PRO:HA	1:C:94:PRO:HD3	1.94	0.40
1:B:82:LEU:HD13	1:B:249:ALA:HB2	2.03	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	537/566 (95%)	505 (94%)	29 (5%)	3 (1%)	30	66
1	B	536/566 (95%)	506 (94%)	26 (5%)	4 (1%)	26	62
1	C	537/566 (95%)	510 (95%)	24 (4%)	3 (1%)	30	66
1	D	537/566 (95%)	506 (94%)	29 (5%)	2 (0%)	39	73
All	All	2147/2264 (95%)	2027 (94%)	108 (5%)	12 (1%)	30	66

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	531	LYS
1	A	539	ILE

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Mol	Chain	Res	Type
1	B	531	LYS
1	D	531	LYS
1	D	230	GLU
1	A	223	CYS
1	B	538	PRO
1	B	188	GLY
1	B	347	SER
1	C	223	CYS
1	C	376	ALA
1	C	531	LYS

### 5.3.2 Protein sidechains [i](#)

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	461/482 (96%)	437 (95%)	24 (5%)	29	63
1	B	460/482 (95%)	436 (95%)	24 (5%)	29	63
1	C	461/482 (96%)	435 (94%)	26 (6%)	26	59
1	D	461/482 (96%)	431 (94%)	30 (6%)	21	51
All	All	1843/1928 (96%)	1739 (94%)	104 (6%)	26	59

All (104) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	THR
1	A	56	ARG
1	A	62	ASP
1	A	100	LYS
1	A	106	LYS
1	A	147	VAL
1	A	159	LEU
1	A	182	LEU
1	A	184	GLN
1	A	207	THR
1	A	230	GLU

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Mol	Chain	Res	Type
1	A	254	LYS
1	A	308	LEU
1	A	321	VAL
1	A	336	LEU
1	A	367	SER
1	A	372	VAL
1	A	380	ARG
1	A	399	THR
1	A	405	VAL
1	A	441	LYS
1	A	494	VAL
1	A	506	SER
1	A	556	SER
1	B	18	GLU
1	B	19	SER
1	B	29	SER
1	B	50	LYS
1	B	53	THR
1	B	79	LYS
1	B	81	LYS
1	B	106	LYS
1	B	132	THR
1	B	137	THR
1	B	148	GLN
1	B	159	LEU
1	B	184	GLN
1	B	254	LYS
1	B	257	THR
1	B	262	ILE
1	B	308	LEU
1	B	311	CYS
1	B	321	VAL
1	B	330	GLN
1	B	368	SER
1	B	377	SER
1	B	380	ARG
1	B	494	VAL
1	C	14	CYS
1	C	21	LEU
1	C	50	LYS
1	C	53	THR
1	C	56	ARG

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Mol	Chain	Res	Type
1	C	100	LYS
1	C	106	LYS
1	C	112	SER
1	C	141	LYS
1	C	158	ARG
1	C	159	LEU
1	C	184	GLN
1	C	196	SER
1	C	207	THR
1	C	222	ARG
1	C	230	GLU
1	C	257	THR
1	C	262	ILE
1	C	308	LEU
1	C	377	SER
1	C	399	THR
1	C	405	VAL
1	C	441	LYS
1	C	453	SER
1	C	492	LEU
1	C	494	VAL
1	D	1	SER
1	D	18	GLU
1	D	51	LYS
1	D	53	THR
1	D	100	LYS
1	D	106	LYS
1	D	112	SER
1	D	130	THR
1	D	158	ARG
1	D	159	LEU
1	D	207	THR
1	D	222	ARG
1	D	226	SER
1	D	230	GLU
1	D	257	THR
1	D	262	ILE
1	D	336	LEU
1	D	355	GLN
1	D	371	SER
1	D	377	SER
1	D	380	ARG

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Mol	Chain	Res	Type
1	D	399	THR
1	D	434	LEU
1	D	441	LYS
1	D	476	SER
1	D	487	SER
1	D	494	VAL
1	D	498	ARG
1	D	508	ARG
1	D	539	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	HIS
1	A	58	GLN
1	A	199	GLN
1	A	562	HIS
1	B	33	HIS
1	B	58	GLN
1	B	199	GLN
1	B	330	GLN
1	C	33	HIS
1	C	58	GLN
1	C	199	GLN
1	C	374	HIS
1	D	33	HIS
1	D	58	GLN
1	D	142	ASN
1	D	199	GLN
1	D	355	GLN
1	D	374	HIS
1	D	502	HIS
1	D	562	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 ⓘ

4 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	3B1	A	601	-	15,16,16	1.41	1 (6%)	15,20,20	1.59	3 (20%)
2	3B1	B	601	-	15,16,16	1.75	3 (20%)	15,20,20	3.05	5 (33%)
2	3B1	C	601	-	15,16,16	1.77	4 (26%)	15,20,20	2.07	5 (33%)
2	3B1	D	601	-	15,16,16	1.47	2 (13%)	15,20,20	1.27	2 (13%)

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	3B1	A	601	-	-	0/8/18/18	0/2/2/2
2	3B1	B	601	-	-	0/8/18/18	0/2/2/2
2	3B1	C	601	-	-	0/8/18/18	0/2/2/2
2	3B1	D	601	-	-	0/8/18/18	0/2/2/2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	3B1	C05-N	-2.14	1.42	1.47
2	B	601	3B1	C08-N01	2.05	1.18	1.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	3B1	C10-C02	2.09	1.43	1.38
2	D	601	3B1	C11-C04	2.69	1.43	1.39
2	B	601	3B1	C11-C04	3.00	1.44	1.39
2	C	601	3B1	C11-C04	3.12	1.44	1.39
2	D	601	3B1	C04-C09	3.19	1.55	1.52
2	A	601	3B1	C08-N01	3.47	1.21	1.14
2	B	601	3B1	C04-C09	4.55	1.57	1.52
2	C	601	3B1	C04-C09	4.66	1.57	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	3B1	C09-C08-N01	-9.35	162.13	177.31
2	C	601	3B1	C09-C08-N01	-4.51	169.98	177.31
2	A	601	3B1	C05-N-C	-2.84	104.08	109.23
2	D	601	3B1	C09-C08-N01	-2.82	172.73	177.31
2	B	601	3B1	C05-N-C	-2.68	104.38	109.23
2	C	601	3B1	C05-N-C	-2.55	104.62	109.23
2	A	601	3B1	C09-C08-N01	-2.13	173.85	177.31
2	B	601	3B1	O-C03-C05	2.21	116.91	111.84
2	D	601	3B1	C03-C05-N	2.34	112.06	109.65
2	C	601	3B1	O-C03-C05	2.58	117.74	111.84
2	B	601	3B1	C03-O-C07	3.27	120.89	109.89
2	C	601	3B1	C07-C-N	3.37	113.11	109.65
2	C	601	3B1	C03-O-C07	3.43	121.46	109.89
2	A	601	3B1	C03-C05-N	3.89	113.64	109.65
2	B	601	3B1	C03-C05-N	4.94	114.72	109.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	3B1	10	0
2	B	601	3B1	2	0
2	C	601	3B1	4	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	545/566 (96%)	-0.34	2 (0%) 93 92	16, 27, 45, 69	0
1	B	544/566 (96%)	-0.31	2 (0%) 93 92	14, 27, 47, 82	0
1	C	545/566 (96%)	-0.30	6 (1%) 82 81	15, 26, 45, 81	0
1	D	545/566 (96%)	-0.32	1 (0%) 95 95	14, 26, 47, 75	0
All	All	2179/2264 (96%)	-0.31	11 (0%) 91 91	14, 26, 46, 82	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	35	ASN	2.8
1	B	15	ALA	2.7
1	A	34	HIS	2.6
1	C	43	ARG	2.6
1	C	14	CYS	2.5
1	A	377	SER	2.5
1	C	10	LEU	2.4
1	C	16	ALA	2.3
1	C	9	ALA	2.2
1	C	15	ALA	2.1
1	B	43	ARG	2.1

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	3B1	A	601	15/15	0.87	0.35	8.69	42,56,66,67	0
2	3B1	C	601	15/15	0.91	0.28	7.63	33,36,39,39	0
2	3B1	D	601	15/15	0.89	0.31	6.53	44,48,52,53	0
2	3B1	B	601	15/15	0.89	0.29	5.17	46,49,53,54	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.