



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:33 PM GMT

PDB ID : 4P73
Title : PheRS in complex with compound 1a
Authors : Ferguson, A.D.
Deposited on : 2014-03-25
Resolution : 3.03 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

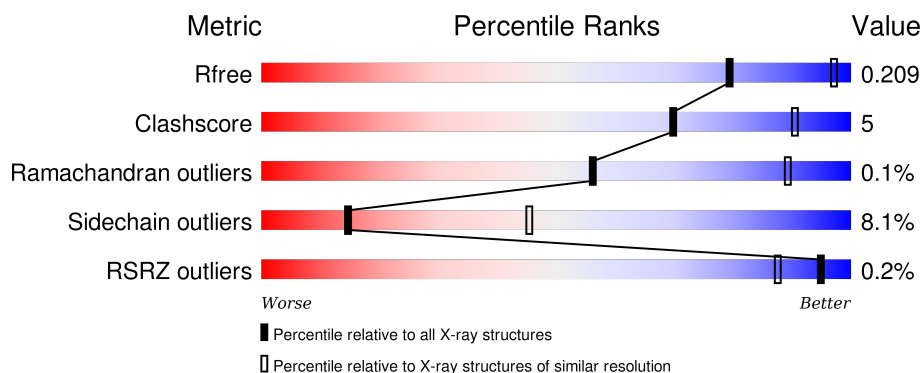
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.03 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	1995 (3.08-3.00)
Clashscore	102246	2351 (3.08-3.00)
Ramachandran outliers	100387	2272 (3.08-3.00)
Sidechain outliers	100360	2275 (3.08-3.00)
RSRZ outliers	91569	2013 (3.08-3.00)

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	792	<div> <div style="width: 83%;"></div> <div style="width: 15%;"></div> <div style="width: 2%;"></div> </div> <div>83% 15% .</div>
1	B	792	<div> <div style="width: 85%;"></div> <div style="width: 13%;"></div> <div style="width: 2%;"></div> </div> <div>85% 13% .</div>
2	C	338	<div> <div style="width: 59%;"></div> <div style="width: 10%;"></div> <div style="width: 29%;"></div> </div> <div>59% 10% . 29%</div>
2	D	338	<div> <div style="width: 57%;"></div> <div style="width: 13%;"></div> <div style="width: 28%;"></div> </div> <div>57% 13% . 28%</div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 16191 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.

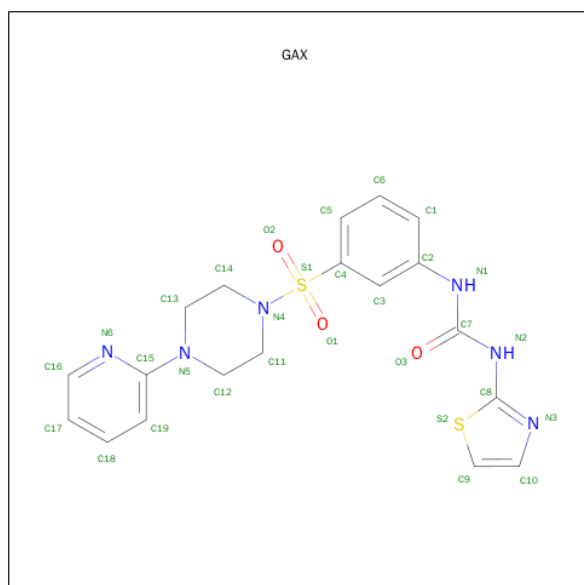
- Molecule 1 is a protein called Phenylalanine-tRNA ligase beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	791	Total	C	N	O	S	0	0	0
			6103	3854	1087	1141	21			
1	B	791	Total	C	N	O	S	0	0	0
			6103	3854	1087	1141	21			

- Molecule 2 is a protein called Phenylalanine-tRNA ligase alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	239	Total	C	N	O	S	0	0	0
			1923	1218	340	351	14			
2	D	243	Total	C	N	O	S	0	0	0
			1952	1238	344	355	15			

- Molecule 3 is 1-{3-[(4-pyridin-2-ylpiperazin-1-yl)sulfonyl]phenyl}-3-(1,3-thiazol-2-yl)urea (three-letter code: GAX) (formula: C₁₉H₂₀N₆O₃S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	C	1	Total	C	N	O	S	0	0
			30	19	6	3	2		
3	D	1	Total	C	N	O	S	0	0
			30	19	6	3	2		

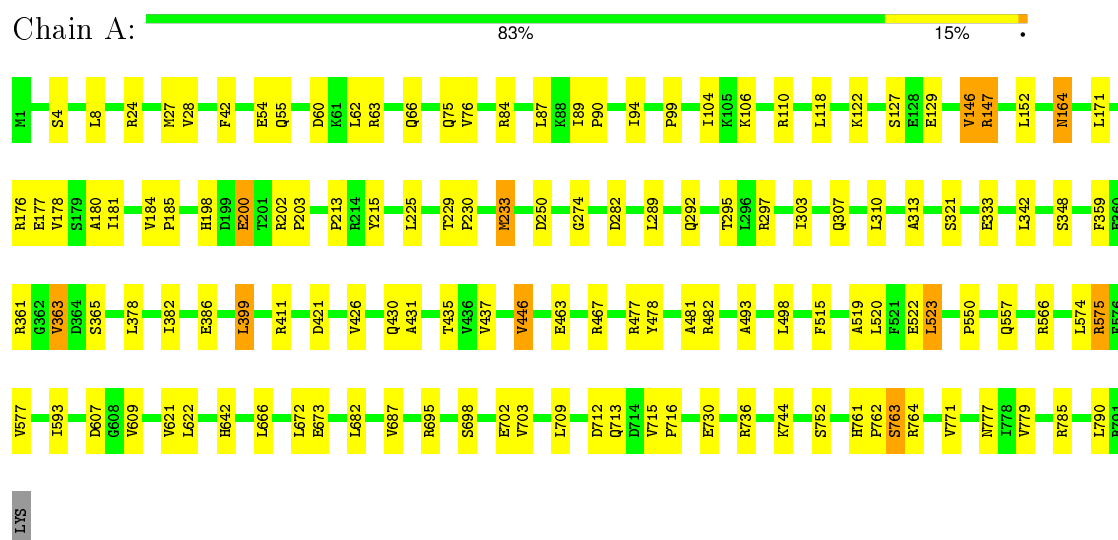
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	20	Total	O	0	0
			20	20		
4	B	22	Total	O	0	0
			22	22		
4	C	4	Total	O	0	0
			4	4		
4	D	4	Total	O	0	0
			4	4		

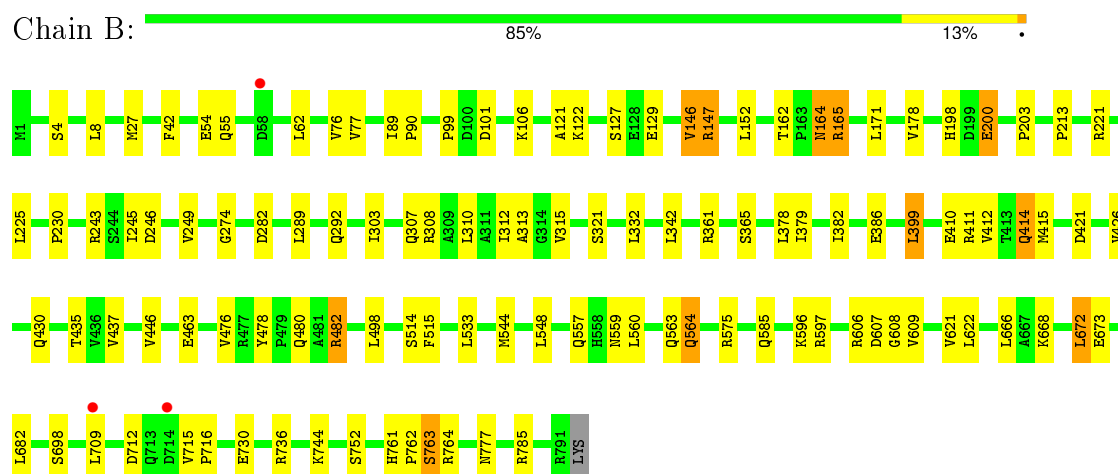
3 Residue-property plots

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: Phenylalanine-tRNA ligase beta subunit

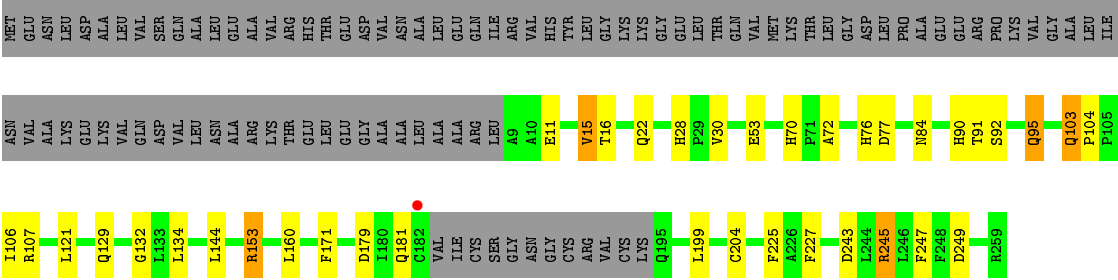


• Molecule 1: Phenylalanine-tRNA ligase beta subunit

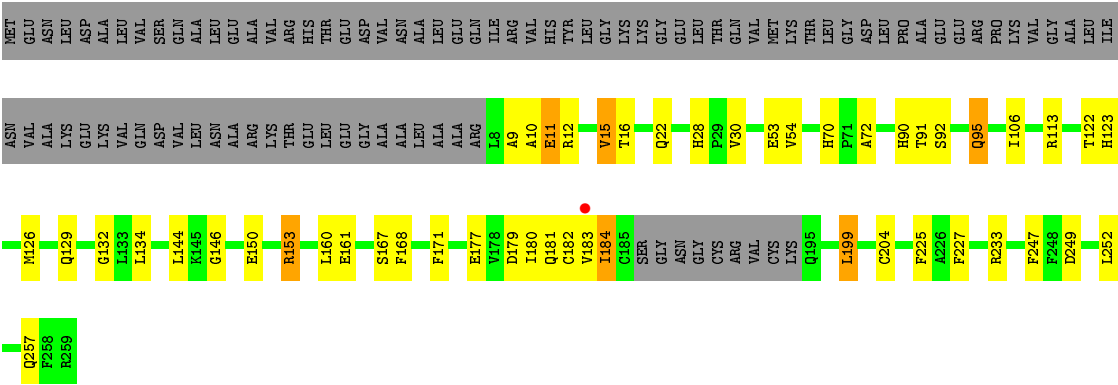


• Molecule 2: Phenylalanine-tRNA ligase alpha subunit





● Molecule 2: Phenylalanine-tRNA ligase alpha subunit



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.29Å 220.05Å 107.93Å 90.00° 101.76° 90.00°	Depositor
Resolution (Å)	49.87 – 3.03 49.87 – 3.03	Depositor EDS
% Data completeness (in resolution range)	99.3 (49.87-3.03) 99.3 (49.87-3.03)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.07 (at 3.01Å)	Xtriage
Refinement program	BUSTER 2.11.5	Depositor
R, R_{free}	0.172 , 0.213 0.171 , 0.209	Depositor DCC
R_{free} test set	2548 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	53.8	Xtriage
Anisotropy	0.501	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 50192 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16191	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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: GAX

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	0.50	1/6221 (0.0%)	0.72	0/8450
1	B	0.50	0/6221	0.72	0/8450
2	C	0.54	0/1972	0.73	3/2664 (0.1%)
2	D	0.53	0/2001	0.73	2/2704 (0.1%)
All	All	0.51	1/16415 (0.0%)	0.72	5/22268 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	233	MET	SD-CE	-5.21	1.48	1.77

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	15	VAL	C-N-CA	6.58	138.16	121.70
2	C	15	VAL	C-N-CA	6.38	137.64	121.70
2	D	153	ARG	CD-NE-CZ	5.77	131.68	123.60
2	C	153	ARG	CD-NE-CZ	5.13	130.78	123.60
2	C	181	GLN	C-N-CA	5.13	134.52	121.70

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	6103	0	6136	62	0
1	B	6103	0	6136	50	0
2	C	1923	0	1846	20	0
2	D	1952	0	1882	32	0
3	C	30	0	20	1	0
3	D	30	0	20	1	0
4	A	20	0	0	0	0
4	B	22	0	0	0	0
4	C	4	0	0	0	0
4	D	4	0	0	0	0
All	All	16191	0	16040	145	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 (145) 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:559:ASN:HD21	2:D:252:LEU:H	1.26	0.82
2:C:15:VAL:HG22	2:C:16:THR:H	1.45	0.81
2:D:15:VAL:HG22	2:D:16:THR:H	1.47	0.80
1:A:566:ARG:HH12	2:C:22:GLN:HE22	1.30	0.78
1:A:695:ARG:H	2:D:257:GLN:HE22	1.35	0.75
1:A:198:HIS:HD2	1:A:200:GLU:H	1.33	0.74
1:A:359:PHE:CD1	1:A:363:VAL:HG21	2.22	0.74
1:A:761:HIS:HD2	1:A:763:SER:H	1.34	0.74
1:B:761:HIS:HD2	1:B:763:SER:H	1.35	0.74
2:C:70:HIS:HD2	2:C:72:ALA:H	1.36	0.73
1:B:198:HIS:HD2	1:B:200:GLU:H	1.34	0.73
2:D:70:HIS:HD2	2:D:72:ALA:H	1.36	0.72
1:A:566:ARG:HH12	2:C:22:GLN:NE2	1.87	0.72
1:B:564:GLN:HG2	2:D:252:LEU:HD12	1.71	0.71
2:C:91:THR:H	2:C:129:GLN:HE21	1.39	0.69
2:D:91:THR:H	2:D:129:GLN:HE21	1.38	0.69
2:C:28:HIS:HD2	2:C:30:VAL:H	1.42	0.67
1:A:478:TYR:O	1:B:482:ARG:HG3	1.95	0.66
2:D:28:HIS:HD2	2:D:30:VAL:H	1.42	0.65
1:B:559:ASN:ND2	2:D:252:LEU:H	1.96	0.64
1:B:414:GLN:HG3	1:B:415:MET:N	2.12	0.63
2:D:167:SER:HB2	2:D:177:GLU:HG3	1.82	0.61
2:D:161:GLU:HB2	2:D:181:GLN:HE21	1.64	0.61
1:A:642:HIS:HE2	2:D:12:ARG:HH21	1.49	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:122:THR:HG23	2:D:123:HIS:CE1	2.37	0.60
1:A:702:GLU:H	1:B:563:GLN:HE22	1.49	0.59
1:B:761:HIS:CD2	1:B:763:SER:H	2.20	0.59
1:A:203:PRO:HG2	1:A:274:GLY:HA3	1.85	0.59
1:A:519:ALA:O	1:A:523:LEU:HD13	2.03	0.59
1:B:761:HIS:CE1	1:B:764:ARG:NH2	2.71	0.58
1:B:164:ASN:H	1:B:164:ASN:HD22	1.51	0.58
1:B:62:LEU:HD22	1:B:77:VAL:HG12	1.85	0.57
1:A:233:MET:CE	1:A:250:ASP:HB3	2.33	0.57
1:B:203:PRO:HG2	1:B:274:GLY:HA3	1.85	0.57
1:A:164:ASN:H	1:A:164:ASN:HD22	1.51	0.57
1:B:761:HIS:CE1	1:B:764:ARG:HH21	2.22	0.56
1:A:477:ARG:HH21	1:B:482:ARG:NH1	2.04	0.56
1:B:564:GLN:HE21	2:D:252:LEU:HB2	1.71	0.56
1:A:593:ILE:HD12	1:A:609:VAL:HG21	1.88	0.56
1:A:761:HIS:CD2	1:A:763:SER:H	2.19	0.55
1:A:712:ASP:O	1:A:715:VAL:HG12	2.07	0.55
1:B:712:ASP:O	1:B:715:VAL:HG12	2.07	0.55
1:B:378:LEU:O	1:B:382:ILE:HG12	2.06	0.55
1:A:359:PHE:CE1	1:A:363:VAL:HG21	2.43	0.54
1:A:378:LEU:O	1:A:382:ILE:HG12	2.06	0.54
1:A:481:ALA:HB2	2:C:107:ARG:NH2	2.23	0.54
1:B:198:HIS:CD2	1:B:200:GLU:HB2	2.44	0.53
1:A:89:ILE:HD13	1:A:118:LEU:HD22	1.91	0.53
1:A:557:GLN:HE21	1:A:672:LEU:HD13	1.73	0.53
1:A:779:VAL:HG11	1:A:790:LEU:HD21	1.90	0.52
2:C:243:ASP:OD1	2:C:245:ARG:HD3	2.09	0.52
1:A:66:GLN:HE22	1:A:75:GLN:HG3	1.75	0.52
1:A:94:ILE:O	1:A:104:ILE:O	2.28	0.52
1:A:28:VAL:HG21	1:A:181:ILE:CD1	2.39	0.51
1:A:198:HIS:CD2	1:A:200:GLU:HB2	2.45	0.51
2:D:132:GLY:HA3	2:D:227:PHE:CZ	2.45	0.51
1:A:229:THR:HG23	1:A:233:MET:HE2	1.92	0.51
1:B:426:VAL:O	1:B:430:GLN:HG3	2.10	0.51
1:A:84:ARG:CZ	1:A:87:LEU:HD21	2.41	0.51
1:A:359:PHE:CD1	1:A:363:VAL:CG2	2.94	0.51
2:C:132:GLY:HA3	2:C:227:PHE:CZ	2.46	0.51
1:A:426:VAL:O	1:A:430:GLN:HG3	2.10	0.50
2:D:91:THR:H	2:D:129:GLN:NE2	2.09	0.50
1:A:8:LEU:HD21	1:A:178:VAL:HG21	1.94	0.50
1:B:698:SER:HB2	1:B:762:PRO:HA	1.93	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:184:ILE:H	2:D:184:ILE:HD12	1.77	0.50
2:C:91:THR:H	2:C:129:GLN:NE2	2.09	0.49
1:B:8:LEU:HD21	1:B:178:VAL:HG21	1.95	0.49
1:A:716:PRO:HA	2:D:15:VAL:HG21	1.93	0.49
1:A:698:SER:HB2	1:A:762:PRO:HA	1.93	0.49
2:C:28:HIS:HE1	2:C:247:PHE:O	1.97	0.48
1:A:213:PRO:HB2	1:A:399:LEU:HD13	1.96	0.48
1:B:121:ALA:HB1	1:B:127:SER:HB3	1.97	0.47
1:A:498:LEU:HD22	1:A:621:VAL:HG13	1.97	0.47
1:A:520:LEU:HA	1:A:523:LEU:HD22	1.96	0.47
1:B:716:PRO:HA	2:C:15:VAL:HG21	1.97	0.47
2:D:179:ASP:HA	2:D:199:LEU:O	2.15	0.47
2:C:144:LEU:HD11	2:C:204:CYS:HB2	1.97	0.47
2:D:144:LEU:HD13	2:D:225:PHE:HB2	1.97	0.47
2:C:76:HIS:HD2	2:C:77:ASP:CG	2.18	0.47
2:C:144:LEU:HD13	2:C:225:PHE:HB2	1.97	0.46
1:B:310:LEU:HD22	1:B:321:SER:HB3	1.98	0.46
1:A:28:VAL:HG21	1:A:181:ILE:HD11	1.98	0.46
1:A:233:MET:HE3	1:A:250:ASP:HB3	1.98	0.46
1:B:312:ILE:HG12	1:B:315:VAL:HG13	1.97	0.46
1:A:566:ARG:NH1	2:C:22:GLN:HE22	2.07	0.46
2:D:144:LEU:HD11	2:D:204:CYS:HB2	1.98	0.46
1:B:213:PRO:HB2	1:B:399:LEU:HD13	1.96	0.46
2:D:28:HIS:HE1	2:D:247:PHE:O	1.99	0.45
1:B:709:LEU:O	1:B:752:SER:HA	2.16	0.45
1:A:180:ALA:HA	1:A:431:ALA:HB1	1.98	0.45
1:A:493:ALA:O	1:A:687:VAL:HG12	2.16	0.45
1:B:42:PHE:HB3	1:B:99:PRO:HD3	1.99	0.45
1:B:585:GLN:NE2	2:D:54:VAL:HB	2.32	0.45
1:A:42:PHE:HB3	1:A:99:PRO:HD3	1.99	0.45
1:B:597:ARG:HA	1:B:609:VAL:CG1	2.47	0.45
1:A:709:LEU:O	1:A:752:SER:HA	2.17	0.45
2:C:179:ASP:HA	2:C:199:LEU:O	2.17	0.44
1:B:557:GLN:HE21	1:B:672:LEU:HD12	1.82	0.44
1:B:498:LEU:HD22	1:B:621:VAL:HG13	1.98	0.44
1:A:90:PRO:HB3	1:A:146:VAL:HG21	1.99	0.44
2:D:180:ILE:HD12	2:D:184:ILE:CD1	2.48	0.44
1:A:303:ILE:HD12	1:A:313:ALA:HB2	2.00	0.44
1:A:520:LEU:HG	1:A:550:PRO:HB3	2.00	0.44
1:B:90:PRO:HB3	1:B:146:VAL:HG21	1.99	0.44
1:B:147:ARG:HA	1:B:152:LEU:HD12	2.00	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:476:VAL:HG13	2:D:146:GLY:HA2	2.00	0.43
1:B:597:ARG:HA	1:B:609:VAL:HG12	2.00	0.43
1:A:229:THR:CG2	1:A:233:MET:HE2	2.48	0.43
1:B:230:PRO:HG3	1:B:382:ILE:HD12	2.01	0.43
3:C:301:GAX:S2	3:C:301:GAX:O3	2.77	0.43
1:B:303:ILE:HD12	1:B:313:ALA:HB2	2.00	0.43
1:A:184:VAL:CG2	1:A:185:PRO:HD2	2.49	0.42
1:A:230:PRO:HG3	1:A:382:ILE:HD12	2.01	0.42
2:C:92:SER:HA	2:C:95:GLN:HG3	2.01	0.42
1:A:297:ARG:NH1	1:A:348:SER:HB2	2.33	0.42
1:B:761:HIS:CD2	1:B:764:ARG:HG2	2.54	0.42
1:A:435:THR:O	1:A:446:VAL:HA	2.20	0.42
2:D:113:ARG:HH11	2:D:126:MET:HE1	1.84	0.42
2:D:92:SER:HA	2:D:95:GLN:HG3	2.02	0.42
1:A:482:ARG:HD3	1:B:478:TYR:O	2.20	0.42
1:B:165:ARG:HD2	1:B:165:ARG:HA	1.83	0.42
2:C:103:GLN:HG2	2:C:104:PRO:HD2	2.01	0.41
1:A:147:ARG:HA	1:A:152:LEU:HD12	2.02	0.41
1:B:198:HIS:HD2	1:B:200:GLU:N	2.11	0.41
1:B:533:LEU:HG	1:B:544:MET:CE	2.50	0.41
1:B:478:TYR:HE1	2:D:150:GLU:HG2	1.85	0.41
2:D:9:ALA:C	2:D:11:GLU:H	2.23	0.41
1:A:703:VAL:HG11	1:A:771:VAL:HG21	2.01	0.41
1:B:435:THR:O	1:B:446:VAL:HA	2.20	0.41
1:A:575:ARG:HD2	1:A:577:VAL:HG23	2.02	0.41
1:A:761:HIS:CD2	1:A:764:ARG:HG2	2.56	0.41
2:D:181:GLN:C	2:D:183:VAL:H	2.23	0.41
1:A:66:GLN:NE2	1:A:75:GLN:HG3	2.34	0.41
3:D:301:GAX:O3	3:D:301:GAX:S2	2.78	0.41
2:C:106:ILE:O	2:C:134:LEU:HA	2.21	0.41
1:B:332:LEU:HD11	1:B:379:ILE:HD12	2.03	0.41
1:A:28:VAL:HG21	1:A:181:ILE:HD13	2.02	0.40
1:A:310:LEU:HD22	1:A:321:SER:HB3	2.02	0.40
2:D:106:ILE:O	2:D:134:LEU:HA	2.21	0.40
1:B:514:SER:OG	2:D:126:MET:HE1	2.20	0.40
1:A:215:TYR:CZ	1:A:333:GLU:HG3	2.57	0.40
1:A:177:GLU:OE2	1:A:467:ARG:NH1	2.55	0.40
1:B:596:LYS:HA	1:B:608:GLY:HA2	2.03	0.40
1:B:162:THR:HB	2:D:168:PHE:HB3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	789/792 (100%)	765 (97%)	24 (3%)	0	100	100
1	B	789/792 (100%)	765 (97%)	24 (3%)	0	100	100
2	C	235/338 (70%)	224 (95%)	11 (5%)	0	100	100
2	D	239/338 (71%)	226 (95%)	11 (5%)	2 (1%)	24	64
All	All	2052/2260 (91%)	1980 (96%)	70 (3%)	2 (0%)	56	89

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	182	CYS
2	D	10	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	645/646 (100%)	590 (92%)	55 (8%)	13	43
1	B	645/646 (100%)	585 (91%)	60 (9%)	11	38
2	C	207/287 (72%)	195 (94%)	12 (6%)	25	62
2	D	211/287 (74%)	199 (94%)	12 (6%)	25	62
All	All	1708/1866 (92%)	1569 (92%)	139 (8%)	15	45

All (139) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	SER
1	A	24	ARG
1	A	27	MET
1	A	54	GLU
1	A	55	GLN
1	A	60	ASP
1	A	62	LEU
1	A	63	ARG
1	A	76	VAL
1	A	106	LYS
1	A	110	ARG
1	A	122	LYS
1	A	127	SER
1	A	129	GLU
1	A	146	VAL
1	A	147	ARG
1	A	164	ASN
1	A	171	LEU
1	A	176	ARG
1	A	200	GLU
1	A	202	ARG
1	A	225	LEU
1	A	282	ASP
1	A	289	LEU
1	A	292	GLN
1	A	295	THR
1	A	307	GLN
1	A	342	LEU
1	A	361	ARG
1	A	363	VAL
1	A	365	SER
1	A	386	GLU
1	A	399	LEU
1	A	411	ARG
1	A	421	ASP
1	A	437	VAL
1	A	446	VAL
1	A	463	GLU
1	A	515	PHE
1	A	522	GLU
1	A	523	LEU
1	A	574	LEU
1	A	575	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	607	ASP
1	A	622	LEU
1	A	666	LEU
1	A	673	GLU
1	A	682	LEU
1	A	713	GLN
1	A	730	GLU
1	A	736	ARG
1	A	744	LYS
1	A	763	SER
1	A	777	ASN
1	A	785	ARG
1	B	4	SER
1	B	27	MET
1	B	54	GLU
1	B	55	GLN
1	B	76	VAL
1	B	89	ILE
1	B	101	ASP
1	B	106	LYS
1	B	122	LYS
1	B	129	GLU
1	B	146	VAL
1	B	147	ARG
1	B	164	ASN
1	B	165	ARG
1	B	171	LEU
1	B	200	GLU
1	B	221	ARG
1	B	225	LEU
1	B	243	ARG
1	B	245	ILE
1	B	246	ASP
1	B	249	VAL
1	B	282	ASP
1	B	289	LEU
1	B	292	GLN
1	B	307	GLN
1	B	308	ARG
1	B	342	LEU
1	B	361	ARG
1	B	365	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	386	GLU
1	B	399	LEU
1	B	410	GLU
1	B	411	ARG
1	B	412	VAL
1	B	414	GLN
1	B	421	ASP
1	B	437	VAL
1	B	463	GLU
1	B	480	GLN
1	B	482	ARG
1	B	515	PHE
1	B	548	LEU
1	B	560	LEU
1	B	564	GLN
1	B	575	ARG
1	B	606	ARG
1	B	607	ASP
1	B	622	LEU
1	B	666	LEU
1	B	668	LYS
1	B	672	LEU
1	B	673	GLU
1	B	682	LEU
1	B	730	GLU
1	B	736	ARG
1	B	744	LYS
1	B	763	SER
1	B	777	ASN
1	B	785	ARG
2	C	11	GLU
2	C	53	GLU
2	C	84	ASN
2	C	90	HIS
2	C	95	GLN
2	C	103	GLN
2	C	121	LEU
2	C	153	ARG
2	C	160	LEU
2	C	171	PHE
2	C	245	ARG
2	C	249	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	D	11	GLU
2	D	22	GLN
2	D	53	GLU
2	D	90	HIS
2	D	95	GLN
2	D	153	ARG
2	D	160	LEU
2	D	171	PHE
2	D	184	ILE
2	D	199	LEU
2	D	233	ARG
2	D	249	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	75	GLN
1	A	164	ASN
1	A	197	GLN
1	A	198	HIS
1	A	261	GLN
1	A	557	GLN
1	A	690	HIS
1	A	713	GLN
1	A	761	HIS
1	B	164	ASN
1	B	198	HIS
1	B	261	GLN
1	B	414	GLN
1	B	557	GLN
1	B	559	ASN
1	B	563	GLN
1	B	564	GLN
1	B	645	GLN
1	B	713	GLN
1	B	761	HIS
2	C	22	GLN
2	C	28	HIS
2	C	70	HIS
2	C	129	GLN
2	D	28	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	D	70	HIS
2	D	76	HIS
2	D	129	GLN
2	D	181	GLN
2	D	223	GLN
2	D	257	GLN

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	$\# Z > 2$
3	GAX	C	301	-	30,33,33	0.51	0	40,46,46	1.03	2 (5%)
3	GAX	D	301	-	30,33,33	0.56	0	40,46,46	0.99	3 (7%)

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
3	GAX	C	301	-	-	0/22/34/34	0/4/4/4
3	GAX	D	301	-	-	0/22/34/34	0/4/4/4

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	C	301	GAX	C8-N2-C7	-4.36	124.80	130.13
3	D	301	GAX	C8-N2-C7	-4.05	125.17	130.13
3	D	301	GAX	C12-N5-C15	2.04	125.06	120.22
3	D	301	GAX	C9-C10-N3	2.69	116.28	109.36
3	C	301	GAX	C9-C10-N3	2.71	116.34	109.36

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
3	C	301	GAX	1	0
3	D	301	GAX	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 [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, 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	791/792 (99%)	-0.36	0 100 100	31, 50, 80, 109	0
1	B	791/792 (99%)	-0.40	3 (0%) 93 80	30, 48, 82, 107	0
2	C	239/338 (70%)	-0.49	1 (0%) 93 80	35, 47, 75, 93	0
2	D	243/338 (71%)	-0.42	1 (0%) 93 80	29, 45, 76, 107	0
All	All	2064/2260 (91%)	-0.40	5 (0%) 95 87	29, 48, 81, 109	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	183	VAL	3.8
1	B	709	LEU	2.4
1	B	714	ASP	2.2
2	C	182	CYS	2.1
1	B	58	ASP	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, 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
3	GAX	C	301	30/30	0.94	0.23	1.64	41,55,71,73	0
3	GAX	D	301	30/30	0.97	0.19	0.90	38,57,68,69	0

6.5 Other polymers [i](#)

There are no such residues in this entry.