



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 04:02 AM UTC

PDB ID : 6PA2 / pdb_00006pa2
Title : E. coli L-asparaginase II mutant (K162M) in complex with L-Asp at pH 5.6
Authors : Lubkowski, J.; Wlodawer, A.
Deposited on : 2019-06-11
Resolution : 1.90 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

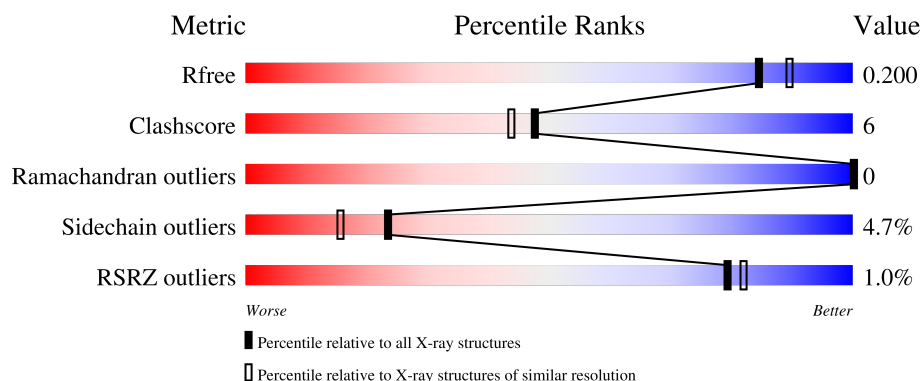
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	334	84% 8% 6%
1	B	334	85% 11% ...
1	C	334	81% 10% 7%
1	D	334	83% 9% 7%

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
3	GOL	B	402	-	X	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10761 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 L-asparaginase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	4	0
			2365	1477	404	475	9			
1	C	311	Total	C	N	O	S	0	2	0
			2345	1467	400	469	9			
1	B	327	Total	C	N	O	S	0	6	0
			2472	1543	421	499	9			
1	D	312	Total	C	N	O	S	0	3	0
			2357	1474	401	472	10			

There are 36 discrepancies between the modelled and reference sequences:

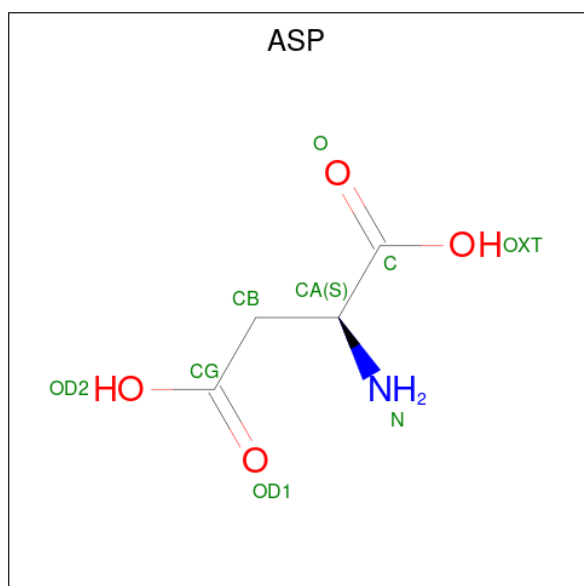
Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	initiating methionine	UNP P00805
A	-6	ASP	-	expression tag	UNP P00805
A	-5	HIS	-	expression tag	UNP P00805
A	-4	HIS	-	expression tag	UNP P00805
A	-3	HIS	-	expression tag	UNP P00805
A	-2	HIS	-	expression tag	UNP P00805
A	-1	HIS	-	expression tag	UNP P00805
A	0	HIS	-	expression tag	UNP P00805
A	162	MET	LYS	engineered mutation	UNP P00805
C	-7	MET	-	initiating methionine	UNP P00805
C	-6	ASP	-	expression tag	UNP P00805
C	-5	HIS	-	expression tag	UNP P00805
C	-4	HIS	-	expression tag	UNP P00805
C	-3	HIS	-	expression tag	UNP P00805
C	-2	HIS	-	expression tag	UNP P00805
C	-1	HIS	-	expression tag	UNP P00805
C	0	HIS	-	expression tag	UNP P00805
C	162	MET	LYS	engineered mutation	UNP P00805
B	-7	MET	-	initiating methionine	UNP P00805
B	-6	ASP	-	expression tag	UNP P00805
B	-5	HIS	-	expression tag	UNP P00805

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	HIS	-	expression tag	UNP P00805
B	-3	HIS	-	expression tag	UNP P00805
B	-2	HIS	-	expression tag	UNP P00805
B	-1	HIS	-	expression tag	UNP P00805
B	0	HIS	-	expression tag	UNP P00805
B	162	MET	LYS	engineered mutation	UNP P00805
D	-7	MET	-	initiating methionine	UNP P00805
D	-6	ASP	-	expression tag	UNP P00805
D	-5	HIS	-	expression tag	UNP P00805
D	-4	HIS	-	expression tag	UNP P00805
D	-3	HIS	-	expression tag	UNP P00805
D	-2	HIS	-	expression tag	UNP P00805
D	-1	HIS	-	expression tag	UNP P00805
D	0	HIS	-	expression tag	UNP P00805
D	162	MET	LYS	engineered mutation	UNP P00805

- Molecule 2 is ASPARTIC ACID (CCD ID: ASP) (formula: $C_4H_7NO_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			9	4	1	4		
2	C	1	Total	C	N	O	0	0
			9	4	1	4		
2	B	1	Total	C	N	O	0	0
			9	4	1	4		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	D	1	Total	C	N	O	0	0
			9	4	1	4		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		

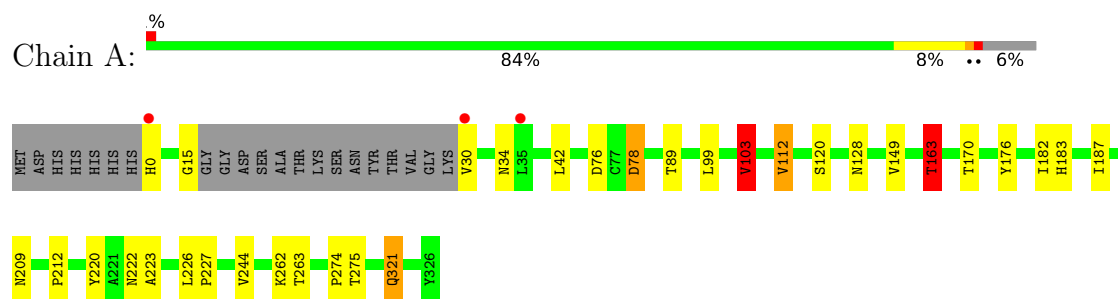
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	324	Total	O	0	2
			326	326		
4	C	277	Total	O	0	5
			282	282		
4	B	287	Total	O	0	0
			287	287		
4	D	282	Total	O	0	3
			285	285		

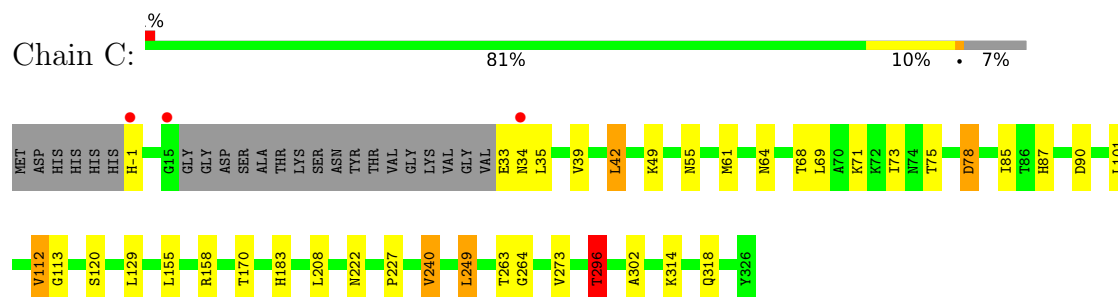
3 Residue-property plots [i](#)

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

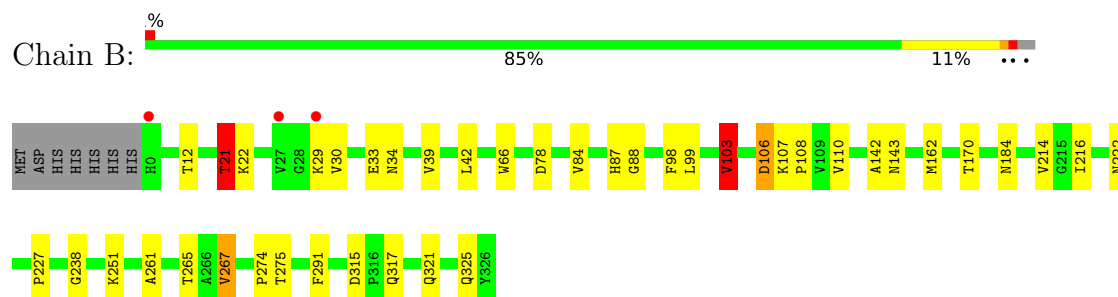
• Molecule 1: L-asparaginase 2



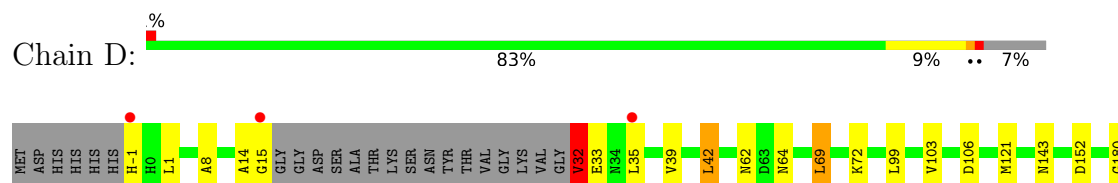
• Molecule 1: L-asparaginase 2



• Molecule 1: L-asparaginase 2



• Molecule 1: L-asparaginase 2





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.18Å 124.47Å 81.74Å 90.00° 112.43° 90.00°	Depositor
Resolution (Å)	26.11 – 1.90 26.11 – 1.90	Depositor EDS
% Data completeness (in resolution range)	91.8 (26.11-1.90) 93.4 (26.11-1.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.13 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.136 , 0.189 0.149 , 0.200	Depositor DCC
R_{free} test set	2494 reflections (2.75%)	wwPDB-VP
Wilson B-factor (Å ²)	22.0	Xtriage
Anisotropy	0.022	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.127 for h,-k,-h-l	Xtriage
Reported twinning fraction	0.154 for H, K, L 0.846 for -H, -K, H+L	Depositor
Outliers	1 of 84439 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10761	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.52% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: GOL

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	1.28	4/2410 (0.2%)	1.21	8/3283 (0.2%)
1	B	1.26	3/2528 (0.1%)	1.19	4/3442 (0.1%)
1	C	1.28	3/2388 (0.1%)	1.22	8/3253 (0.2%)
1	D	1.25	5/2406 (0.2%)	1.22	8/3276 (0.2%)
All	All	1.27	15/9732 (0.2%)	1.21	28/13254 (0.2%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	62	ASN	C-O	6.29	1.31	1.23
1	A	163	THR	CA-C	6.05	1.60	1.52
1	B	216	ILE	CA-C	5.98	1.60	1.52
1	C	273	VAL	CA-C	5.65	1.58	1.53
1	A	212	PRO	N-CA	-5.59	1.40	1.47
1	D	180	GLY	C-O	5.58	1.29	1.23
1	A	128	ASN	CA-C	5.55	1.59	1.52
1	A	275	THR	C-O	5.40	1.30	1.23
1	C	155	LEU	C-O	-5.36	1.17	1.24
1	D	296	THR	CB-CG2	-5.24	1.35	1.52
1	D	273	VAL	CA-C	5.19	1.57	1.53
1	C	113	GLY	N-CA	5.14	1.52	1.45
1	B	251	LYS	N-CA	5.13	1.52	1.46
1	B	275	THR	CA-C	-5.07	1.46	1.52
1	D	223	ALA	C-O	-5.04	1.17	1.23

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	272	ARG	NE-CZ-NH1	9.88	131.38	121.50
1	D	272	ARG	NE-CZ-NH2	-8.57	111.49	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	112	VAL	CB-CA-C	-8.17	96.08	110.71
1	C	112	VAL	CB-CA-C	-7.98	96.81	111.18
1	A	163	THR	N-CA-CB	-7.59	100.49	110.59
1	D	272	ARG	CB-CG-CD	7.38	128.27	111.30
1	C	240	VAL	N-CA-CB	7.29	119.74	111.21
1	C	170	THR	N-CA-C	6.58	119.29	111.33
1	A	274	PRO	N-CA-C	6.33	121.77	113.86
1	A	78	ASP	N-CA-CB	6.21	119.48	110.47
1	A	103	VAL	N-CA-CB	6.12	118.64	110.26
1	B	274	PRO	N-CA-C	6.01	121.37	113.86
1	A	321	GLN	N-CA-CB	5.89	118.87	110.16
1	D	32	VAL	CB-CA-C	-5.81	100.36	111.40
1	C	264	GLY	N-CA-C	5.72	123.70	114.90
1	C	296	THR	N-CA-CB	-5.67	100.46	110.39
1	B	21	THR	N-CA-CB	-5.59	103.17	110.88
1	D	281	ASP	N-CA-C	5.56	119.25	111.74
1	B	103	VAL	N-CA-CB	5.52	118.74	110.13
1	D	272	ARG	CD-NE-CZ	5.52	132.13	124.40
1	A	170	THR	N-CA-C	5.43	117.90	111.33
1	D	103	VAL	N-CA-C	-5.29	101.91	108.89
1	B	170	THR	N-CA-C	5.21	116.96	111.28
1	A	321	GLN	CA-CB-CG	5.15	124.39	114.10
1	C	302	ALA	N-CA-C	-5.15	105.85	111.82
1	C	120	SER	N-CA-C	5.06	117.06	110.53
1	C	112	VAL	CG1-CB-CG2	5.05	121.92	110.80
1	D	264	GLY	N-CA-C	5.03	120.78	114.25

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2365	0	2356	18	0
1	B	2472	0	2468	44	0
1	C	2345	0	2338	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2357	0	2353	24	0
2	A	9	0	3	0	0
2	B	9	0	3	1	0
2	C	9	0	3	0	0
2	D	9	0	3	0	0
3	B	6	0	7	26	0
4	A	326	0	0	13	1
4	B	287	0	0	11	1
4	C	282	0	0	8	0
4	D	285	0	0	16	0
All	All	10761	0	9534	107	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 (107) 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:106:ASP:C	3:B:402:GOL:O2	1.89	1.15
1:B:143[A]:ASN:ND2	3:B:402:GOL:O1	2.01	0.93
1:B:22:LYS:HD3	4:B:522:HOH:O	1.78	0.82
1:B:143[A]:ASN:ND2	3:B:402:GOL:C1	2.44	0.80
1:D:8:ALA:HB2	1:D:32:VAL:HG13	1.63	0.79
1:C:-1:HIS:HA	4:C:732:HOH:O	1.83	0.79
1:B:143[A]:ASN:H	3:B:402:GOL:C1	1.95	0.79
1:D:188:ASP:O	4:D:501:HOH:O	2.01	0.78
1:A:176:TYR:OH	4:D:501:HOH:O	2.02	0.77
1:B:107:LYS:N	3:B:402:GOL:O2	2.17	0.76
1:B:143[A]:ASN:CG	3:B:402:GOL:H11	2.11	0.76
1:D:184:ASN:HB2	4:D:749[B]:HOH:O	1.86	0.74
1:C:296:THR:HG23	4:C:511:HOH:O	1.90	0.72
1:C:314:LYS:HE3	4:C:678:HOH:O	1.90	0.71
1:C:183:HIS:HD2	4:B:749:HOH:O	1.72	0.71
1:A:120:SER:OG	4:A:501:HOH:O	2.09	0.70
1:C:33:GLU:N	4:C:501:HOH:O	2.24	0.70
1:B:106:ASP:CG	3:B:402:GOL:HO1	2.00	0.69
1:A:209[B]:ASN:OD1	4:A:502:HOH:O	2.09	0.69
1:B:87:HIS:HD2	1:B:88:GLY:O	1.74	0.68
1:A:149:VAL:HG12	4:A:574:HOH:O	1.93	0.68
1:B:321[A]:GLN:NE2	4:B:501:HOH:O	2.26	0.68
1:D:106:ASP:OD1	4:D:502:HOH:O	2.13	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:296:THR:CG2	4:C:511:HOH:O	2.42	0.66
1:B:33[B]:GLU:HA	1:B:33[B]:GLU:OE1	1.97	0.65
1:B:143[A]:ASN:ND2	3:B:402:GOL:HO1	1.94	0.65
1:A:15:GLY:C	4:A:528:HOH:O	2.39	0.64
1:C:39:VAL:HB	1:C:42:LEU:HD22	1.79	0.64
1:D:199[A]:SER:OG	4:D:503:HOH:O	2.14	0.63
1:B:162:MET:HE1	4:B:518:HOH:O	1.98	0.63
4:B:663:HOH:O	1:D:272:ARG:HD2	1.99	0.62
1:B:106:ASP:HB3	3:B:402:GOL:O1	1.99	0.62
1:B:33[A]:GLU:HG2	4:B:506:HOH:O	1.98	0.62
1:B:143[A]:ASN:H	3:B:402:GOL:C2	2.14	0.61
1:D:143:ASN:CG	4:D:513:HOH:O	2.44	0.61
3:B:402:GOL:H32	4:B:523:HOH:O	2.01	0.61
1:B:321[B]:GLN:HG2	1:B:325:GLN:HE21	1.66	0.60
1:B:106:ASP:O	3:B:402:GOL:O2	2.19	0.60
1:B:33[A]:GLU:HG3	4:B:780:HOH:O	2.01	0.60
1:B:33[A]:GLU:CG	4:B:506:HOH:O	2.51	0.58
1:D:15:GLY:C	4:D:593:HOH:O	2.45	0.58
4:A:513:HOH:O	1:B:21:THR:HG21	2.03	0.58
1:B:106:ASP:CB	3:B:402:GOL:O1	2.52	0.58
1:B:321[B]:GLN:HG2	1:B:325:GLN:NE2	2.18	0.58
1:D:296:THR:HG23	4:D:522:HOH:O	2.03	0.58
1:B:143[A]:ASN:CG	3:B:402:GOL:C1	2.76	0.57
1:D:296:THR:CG2	4:D:522:HOH:O	2.53	0.57
1:A:183:HIS:HE1	4:A:781:HOH:O	1.87	0.56
1:D:39:VAL:HB	1:D:42:LEU:HD22	1.88	0.56
1:C:249:LEU:N	1:C:249:LEU:HD22	2.21	0.56
1:A:99:LEU:O	1:A:103:VAL:HG13	2.06	0.55
1:C:64:ASN:O	1:C:68[A]:THR:HG23	2.05	0.55
1:B:99:LEU:O	1:B:103:VAL:HG13	2.08	0.54
1:B:315:ASP:OD1	1:B:317[A]:GLN:NE2	2.40	0.54
1:C:73:ILE:HD11	1:C:85:ILE:HD11	1.90	0.54
1:D:251:LYS:HB2	4:D:746:HOH:O	2.08	0.53
1:D:72:LYS:HE2	4:D:527:HOH:O	2.07	0.52
1:C:314:LYS:CE	4:C:678:HOH:O	2.54	0.51
1:D:14:ALA:HA	1:D:35:LEU:HD21	1.91	0.51
1:A:76:ASP:HB3	4:A:716:HOH:O	2.10	0.51
1:A:183:HIS:HD2	4:A:786:HOH:O	1.94	0.51
1:B:12:THR:OG1	2:B:401:ASP:CG	2.53	0.51
1:A:263:THR:CG2	4:A:504:HOH:O	2.59	0.50
1:D:64:ASN:ND2	4:D:510:HOH:O	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:227:PRO:HB3	1:D:227:PRO:HB3	1.92	0.50
1:B:106:ASP:CA	3:B:402:GOL:O2	2.59	0.50
1:B:142:ALA:HA	3:B:402:GOL:H2	1.92	0.49
1:A:89:THR:OG1	4:A:503:HOH:O	2.20	0.49
1:B:143[A]:ASN:H	3:B:402:GOL:C3	2.25	0.48
1:B:184:ASN:HB2	4:B:694:HOH:O	2.13	0.48
1:D:106:ASP:OD1	1:D:106:ASP:N	2.37	0.48
1:B:84:VAL:HA	1:B:110:VAL:O	2.14	0.47
1:B:143[A]:ASN:H	3:B:402:GOL:H31	1.77	0.47
1:C:78:ASP:OD1	1:C:78:ASP:N	2.48	0.47
1:D:32:VAL:N	4:D:515:HOH:O	2.47	0.47
1:A:321:GLN:NE2	4:A:520:HOH:O	2.48	0.46
1:C:39:VAL:HG13	1:D:121[B]:MET:HE1	1.98	0.46
1:B:66:TRP:HB3	1:B:98:PHE:CE2	2.51	0.46
1:B:107:LYS:CA	3:B:402:GOL:O2	2.64	0.46
1:C:71:LYS:O	1:C:75:THR:HG23	2.15	0.46
1:A:163:THR:HG21	4:A:725:HOH:O	2.16	0.45
1:C:33:GLU:N	4:C:518:HOH:O	2.48	0.45
1:B:321[B]:GLN:CG	1:B:325:GLN:HE21	2.30	0.44
1:A:0:HIS:CE1	4:A:511:HOH:O	2.69	0.44
1:C:39:VAL:HG13	1:D:121[B]:MET:CE	2.48	0.44
1:D:35:LEU:CD2	4:D:703:HOH:O	2.65	0.44
1:A:227:PRO:HB3	1:C:227:PRO:HB3	2.01	0.43
1:B:214:VAL:HA	1:B:238:GLY:O	2.18	0.43
3:B:402:GOL:H31	4:B:670:HOH:O	2.19	0.43
1:C:263:THR:HG22	1:C:263:THR:O	2.18	0.43
1:C:61:MET:HE3	1:C:87:HIS:NE2	2.34	0.43
1:C:158:ARG:HH12	1:C:296:THR:HG22	1.84	0.42
1:A:182:ILE:HG12	1:A:187:ILE:HG12	2.00	0.42
1:B:106:ASP:CB	3:B:402:GOL:HO1	2.33	0.42
1:B:106:ASP:HA	3:B:402:GOL:O2	2.18	0.42
1:C:68[A]:THR:HG22	4:C:677:HOH:O	2.19	0.42
1:A:220:TYR:CE2	1:A:223:ALA:HA	2.55	0.41
1:B:261:ALA:HA	1:B:265:THR:O	2.21	0.41
1:B:108:PRO:HA	3:B:402:GOL:O3	2.21	0.41
1:B:142:ALA:CB	3:B:402:GOL:H2	2.51	0.41
1:A:244:VAL:HB	1:C:90:ASP:HB3	2.03	0.41
1:D:69:LEU:HB3	1:D:99:LEU:HD21	2.02	0.41
1:D:325:GLN:NE2	4:D:531:HOH:O	2.54	0.41
1:D:35:LEU:HD23	4:D:703:HOH:O	2.21	0.41
1:B:267:VAL:HG22	1:B:291:PHE:CD1	2.57	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)
4:A:692:HOH:O	4:B:502:HOH:O[1_554]	2.08	0.12

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	313/334 (94%)	307 (98%)	6 (2%)	0	100	100
1	B	331/334 (99%)	321 (97%)	10 (3%)	0	100	100
1	C	309/334 (92%)	304 (98%)	5 (2%)	0	100	100
1	D	312/334 (93%)	304 (97%)	8 (3%)	0	100	100
All	All	1265/1336 (95%)	1236 (98%)	29 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	261/274 (95%)	251 (96%)	10 (4%)	29	21
1	B	273/274 (100%)	262 (96%)	11 (4%)	28	20
1	C	258/274 (94%)	242 (94%)	16 (6%)	16	9
1	D	261/274 (95%)	248 (95%)	13 (5%)	22	14

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1053/1096 (96%)	1003 (95%)	50 (5%)	23	15

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	VAL
1	A	34	ASN
1	A	42	LEU
1	A	78	ASP
1	A	103	VAL
1	A	112	VAL
1	A	163	THR
1	A	222	ASN
1	A	226	LEU
1	A	262	LYS
1	C	34	ASN
1	C	35	LEU
1	C	42	LEU
1	C	49	LYS
1	C	55	ASN
1	C	69	LEU
1	C	78	ASP
1	C	101	LEU
1	C	112	VAL
1	C	129	LEU
1	C	208	LEU
1	C	222	ASN
1	C	240	VAL
1	C	249	LEU
1	C	296	THR
1	C	318	GLN
1	B	21	THR
1	B	29	LYS
1	B	30	VAL
1	B	34	ASN
1	B	39	VAL
1	B	42	LEU
1	B	78	ASP
1	B	103	VAL
1	B	106	ASP
1	B	222	ASN
1	B	267	VAL

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Mol	Chain	Res	Type
1	D	-1	HIS
1	D	1	LEU
1	D	32	VAL
1	D	33	GLU
1	D	42	LEU
1	D	69	LEU
1	D	152	ASP
1	D	199[A]	SER
1	D	199[B]	SER
1	D	222	ASN
1	D	226	LEU
1	D	257	LEU
1	D	296	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	183	HIS
1	A	248	ASN
1	A	280	GLN
1	A	321	GLN
1	C	-1	HIS
1	C	55	ASN
1	C	183	HIS
1	C	248	ASN
1	C	317	GLN
1	C	324	ASN
1	B	87	HIS
1	B	219	ASN
1	B	325	GLN
1	D	248	ASN
1	D	318	GLN
1	D	321	GLN
1	D	324	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	ASP	A	401	-	7,8,8	0.90	0	6,10,10	1.10	0
2	ASP	D	401	-	7,8,8	1.69	2 (28%)	6,10,10	1.34	1 (16%)
3	GOL	B	402	-	5,5,5	1.39	1 (20%)	5,5,5	1.48	1 (20%)
2	ASP	C	401	-	7,8,8	1.88	3 (42%)	6,10,10	1.95	1 (16%)
2	ASP	B	401	-	7,8,8	1.99	2 (28%)	6,10,10	1.37	1 (16%)

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	ASP	A	401	-	-	2/8/8/8	-
2	ASP	D	401	-	-	4/8/8/8	-
3	GOL	B	402	-	-	4/4/4/4	-
2	ASP	C	401	-	-	4/8/8/8	-
2	ASP	B	401	-	-	2/8/8/8	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	ASP	OD1-CG	3.92	1.34	1.22
2	C	401	ASP	OXT-C	-3.03	1.21	1.30
3	B	402	GOL	O2-C2	-2.98	1.34	1.43
2	D	401	ASP	OD1-CG	2.80	1.31	1.22
2	C	401	ASP	OD1-CG	2.73	1.31	1.22
2	D	401	ASP	OXT-C	-2.28	1.23	1.30
2	B	401	ASP	OD2-CG	-2.14	1.23	1.30
2	C	401	ASP	OD2-CG	2.10	1.37	1.30

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	ASP	OXT-C-O	-4.09	114.81	124.08
3	B	402	GOL	O1-C1-C2	3.00	123.86	110.38
2	B	401	ASP	OXT-C-O	-2.76	117.83	124.08
2	D	401	ASP	OD2-CG-CB	2.21	120.87	114.00

There are no chirality outliers.

All (16) torsion outliers are listed below:

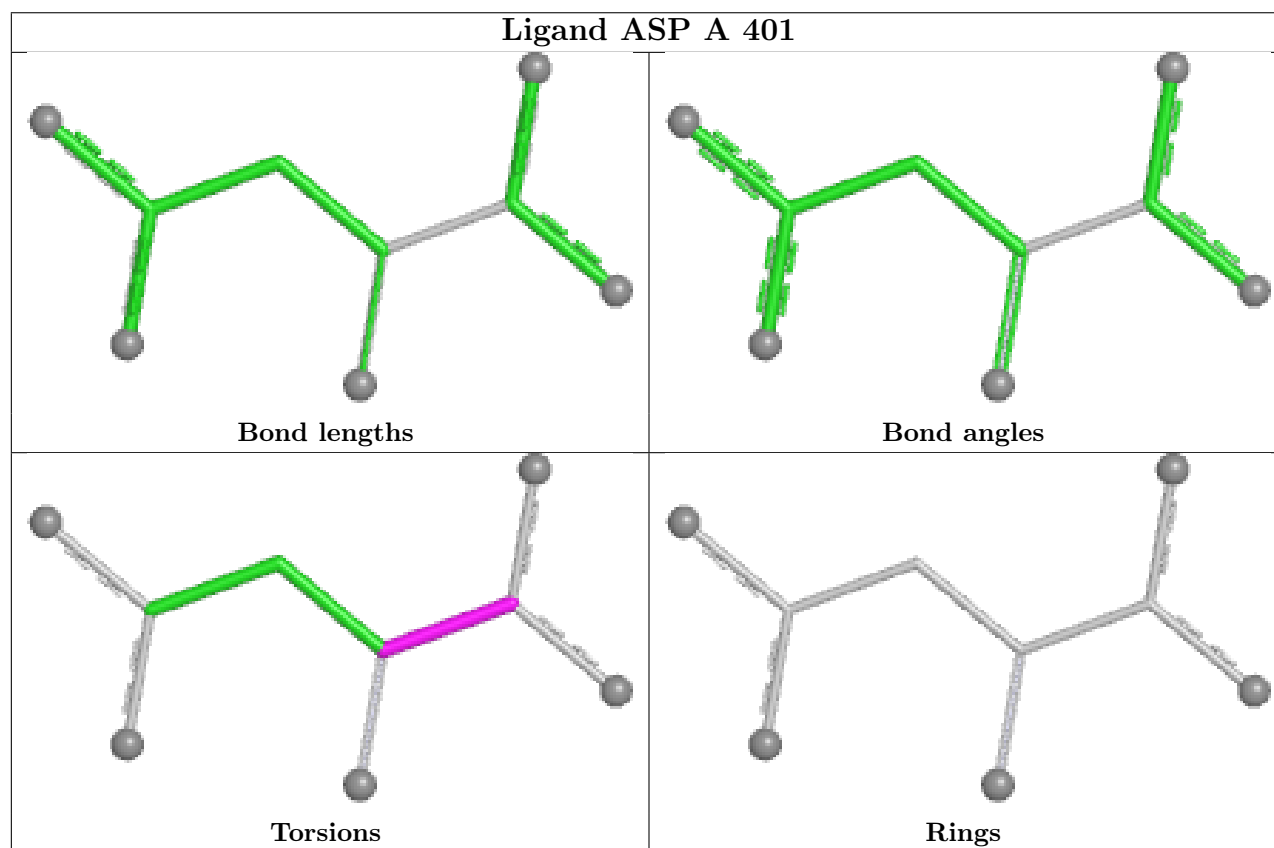
Mol	Chain	Res	Type	Atoms
2	A	401	ASP	O-C-CA-N
2	C	401	ASP	O-C-CA-N
2	B	401	ASP	O-C-CA-N
2	D	401	ASP	O-C-CA-N
3	B	402	GOL	O1-C1-C2-C3
2	C	401	ASP	OXT-C-CA-N
2	D	401	ASP	OXT-C-CA-N
2	B	401	ASP	OXT-C-CA-N
2	A	401	ASP	OXT-C-CA-N
3	B	402	GOL	O2-C2-C3-O3
3	B	402	GOL	O1-C1-C2-O2
2	C	401	ASP	OXT-C-CA-CB
2	C	401	ASP	O-C-CA-CB
2	D	401	ASP	O-C-CA-CB
2	D	401	ASP	OXT-C-CA-CB
3	B	402	GOL	C1-C2-C3-O3

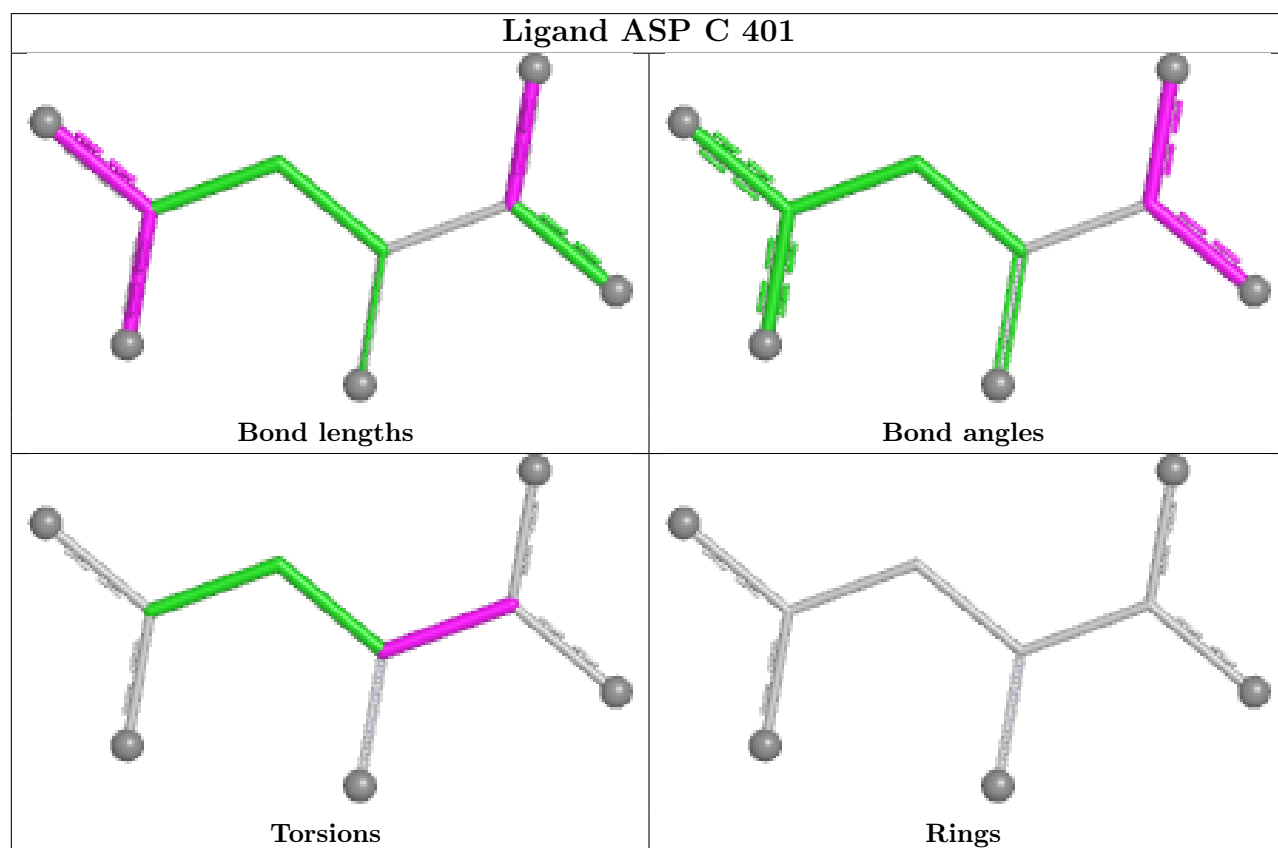
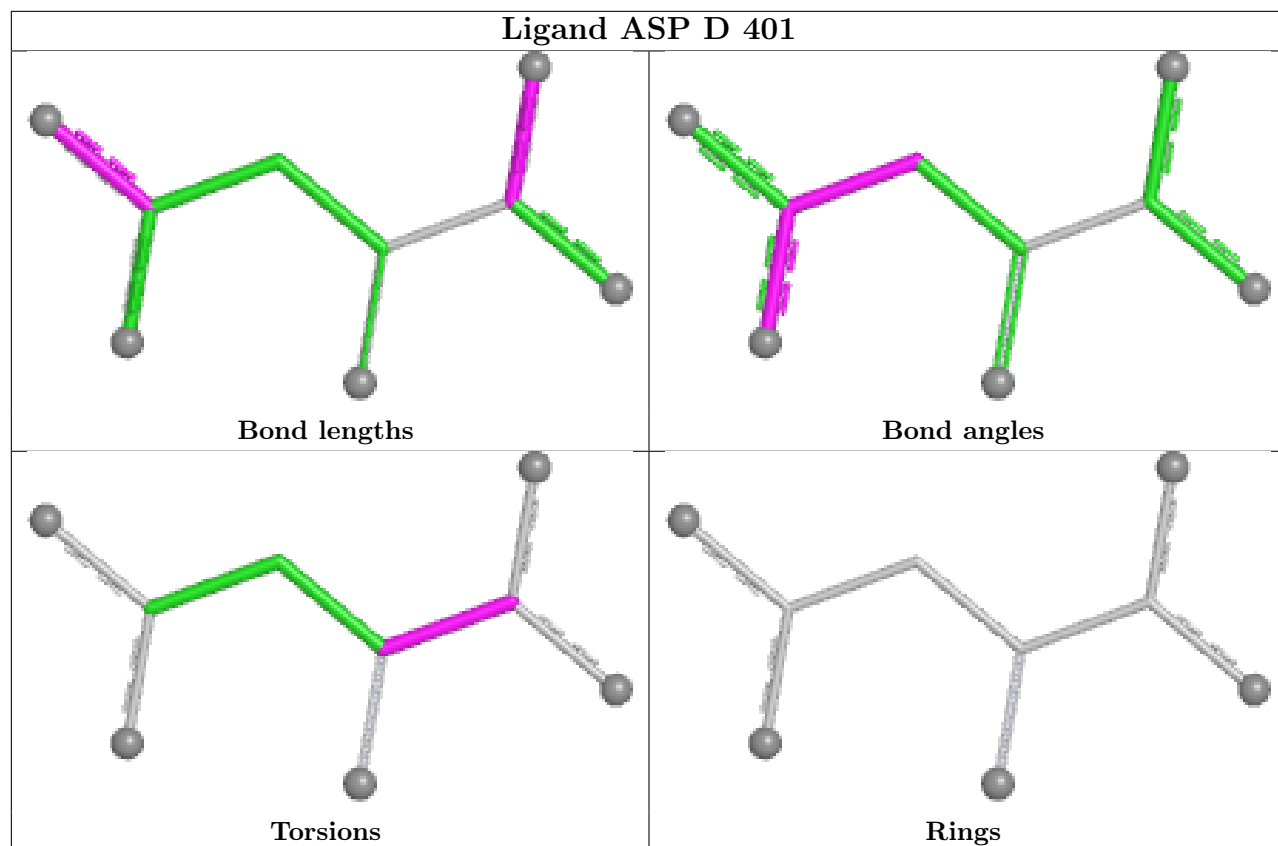
There are no ring outliers.

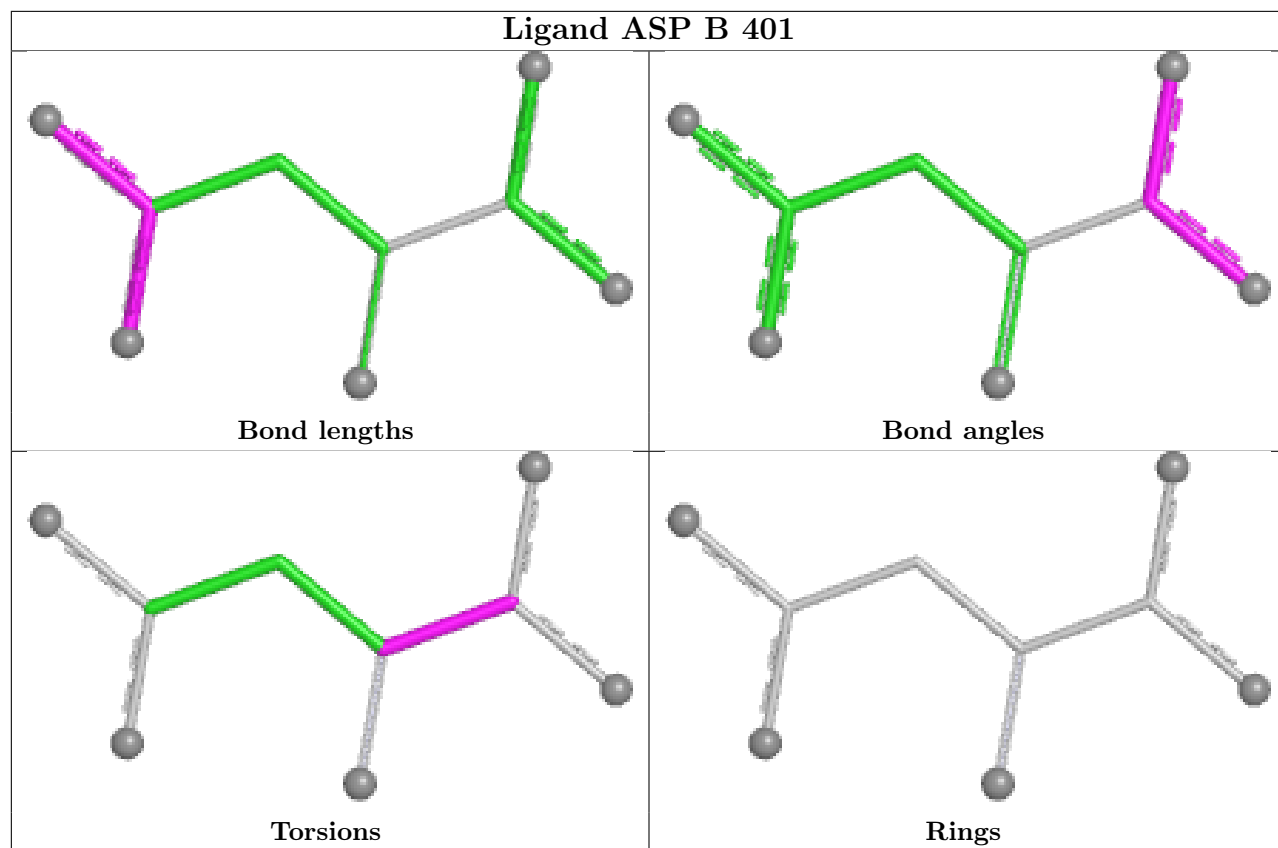
2 monomers are involved in 27 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	GOL	26	0
2	B	401	ASP	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







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	313/334 (93%)	-0.31	3 (0%) 79 82	13, 24, 39, 62	4 (1%)
1	B	327/334 (97%)	-0.16	3 (0%) 81 83	17, 26, 42, 80	6 (1%)
1	C	311/334 (93%)	-0.20	3 (0%) 79 82	17, 25, 40, 68	2 (0%)
1	D	312/334 (93%)	-0.24	3 (0%) 79 82	16, 24, 41, 72	3 (0%)
All	All	1263/1336 (94%)	-0.23	12 (0%) 79 82	13, 25, 41, 80	15 (1%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	15	GLY	3.9
1	B	29	LYS	2.8
1	A	35	LEU	2.8
1	A	0	HIS	2.7
1	A	30	VAL	2.7
1	C	-1	HIS	2.6
1	C	15	GLY	2.5
1	D	-1	HIS	2.4
1	D	35	LEU	2.4
1	B	0	HIS	2.2
1	C	34	ASN	2.2
1	B	27	VAL	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

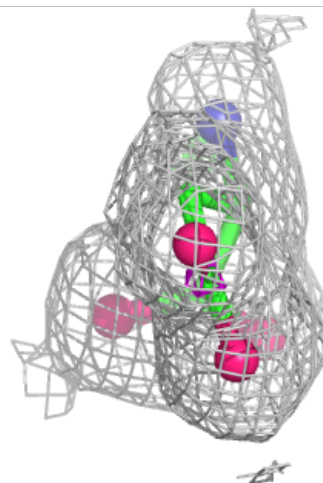
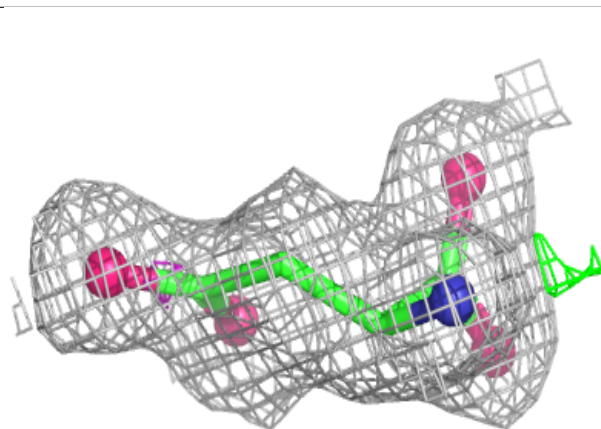
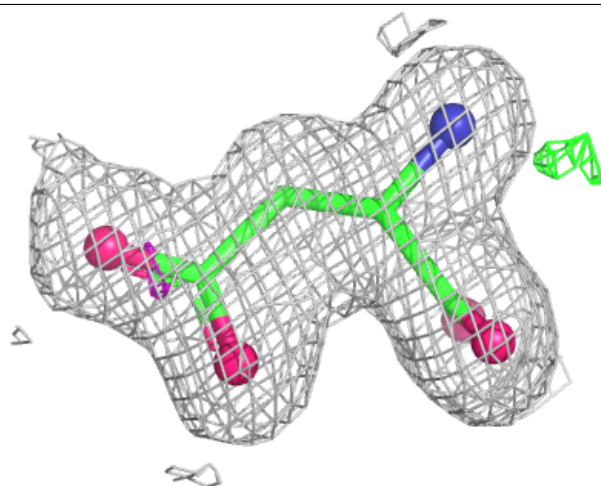
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	B	402	6/6	0.73	0.18	30,30,31,34	6
2	ASP	C	401	9/9	0.93	0.08	22,24,28,28	0
2	ASP	D	401	9/9	0.96	0.06	20,23,27,33	0
2	ASP	B	401	9/9	0.96	0.07	23,24,28,29	0
2	ASP	A	401	9/9	0.97	0.06	20,23,25,32	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

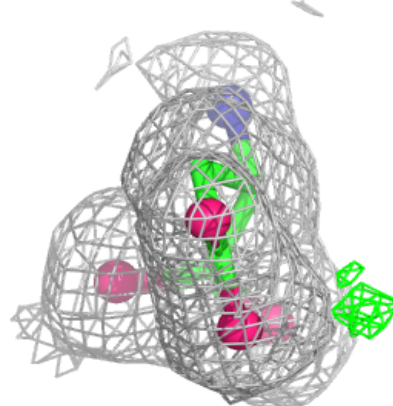
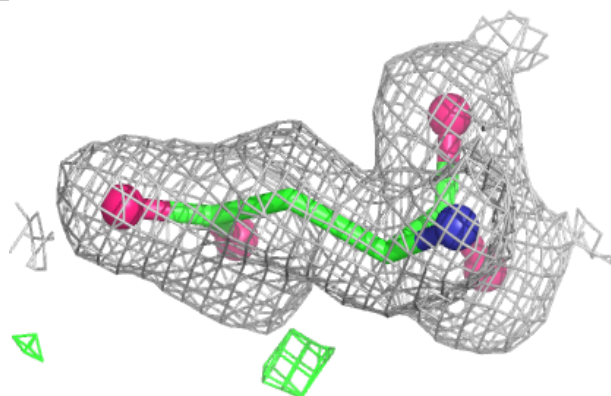
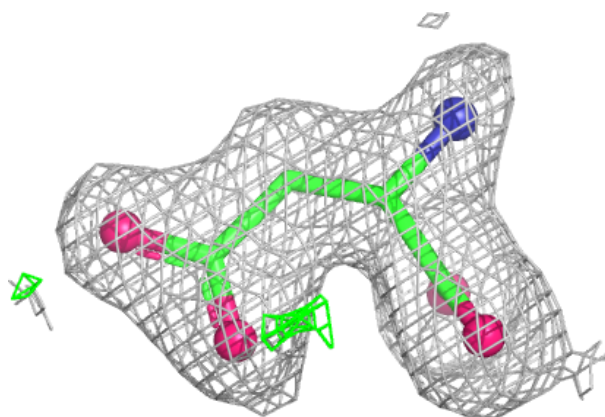
Electron density around ASP C 401:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



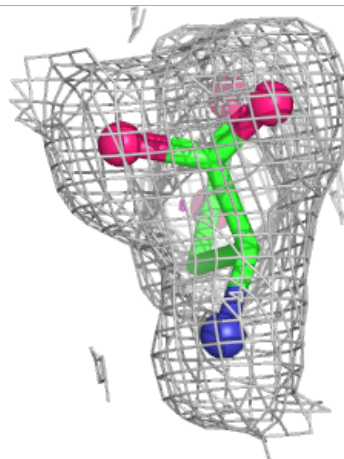
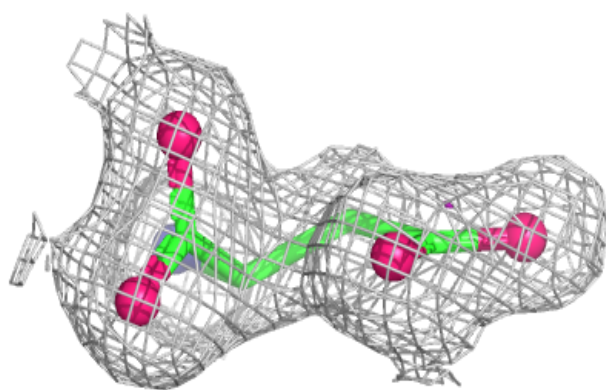
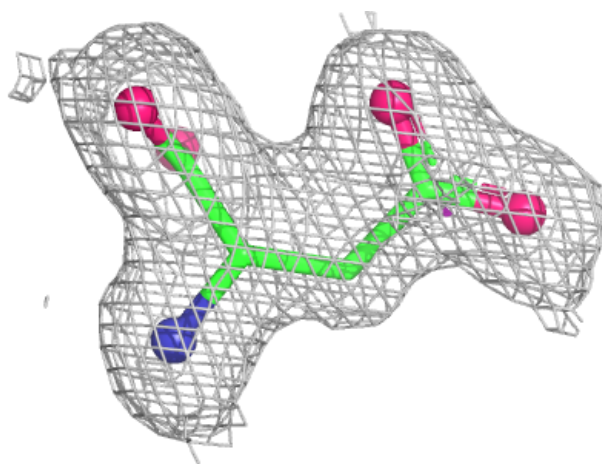
Electron density around ASP D 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



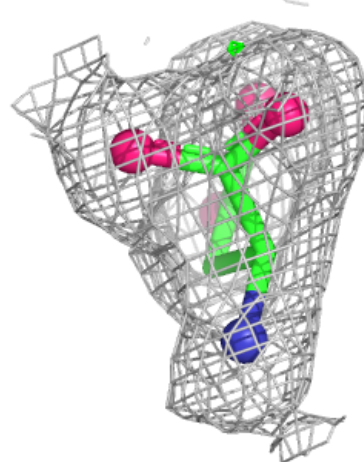
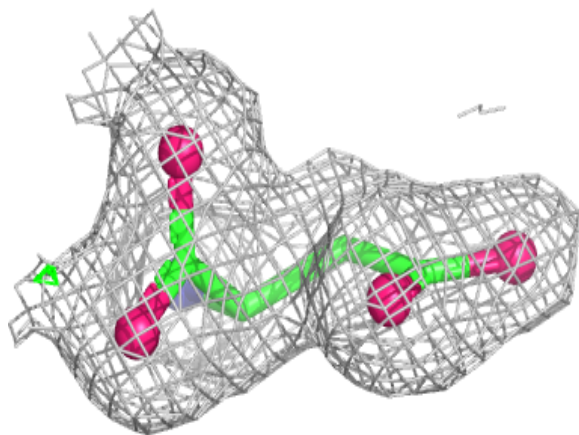
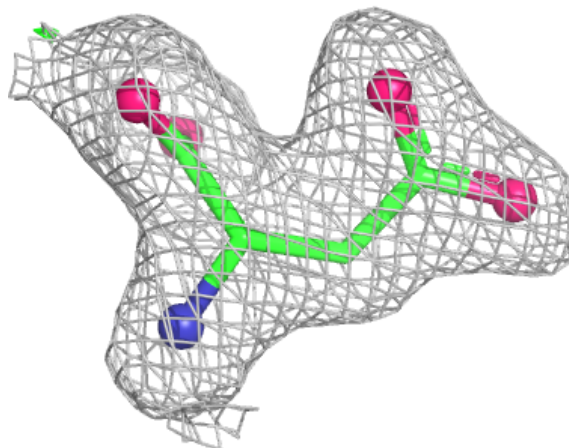
Electron density around ASP B 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ASP A 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.