



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:07 AM GMT

PDB ID : 2W03
Title : CO-COMPLEX STRUCTURE OF ACHROMOBACTIN SYNTHETASE
PROTEIN D (ACSD) WITH ADENOSINE, SULFATE AND CITRATE
FROM PECTOBACTERIUM CHRYSANTHEMI
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Deposited on : 2008-08-08
Resolution : 2.95 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

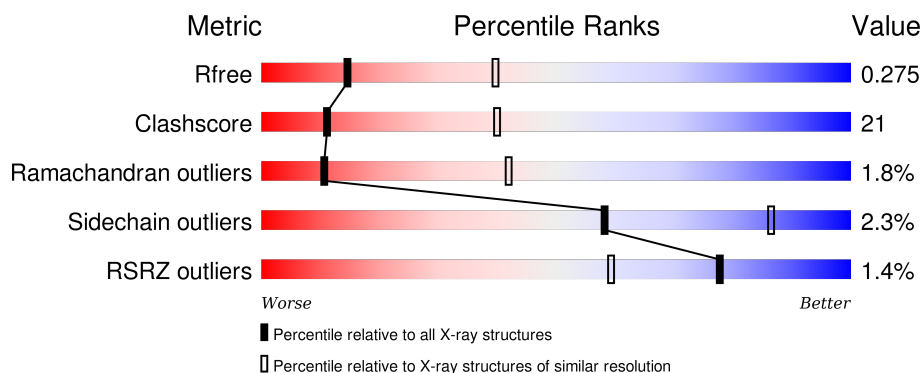
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	2184 (3.00-2.92)
Clashscore	102246	2552 (3.00-2.92)
Ramachandran outliers	100387	2468 (3.00-2.92)
Sidechain outliers	100360	2471 (3.00-2.92)
RSRZ outliers	91569	2201 (3.00-2.92)

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	620	<div> <div> <div></div> <div>55%</div> <div>35%</div> <div>7%</div> </div> <div> <div>2%</div> <div>52%</div> <div>39%</div> <div>7%</div> </div> </div>
1	B	620	<div> <div> <div></div> <div>55%</div> <div>35%</div> <div>7%</div> </div> <div> <div>2%</div> <div>52%</div> <div>39%</div> <div>7%</div> </div> </div>

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	SO4	B	1589	-	-	X	-

2 Entry composition [i](#)

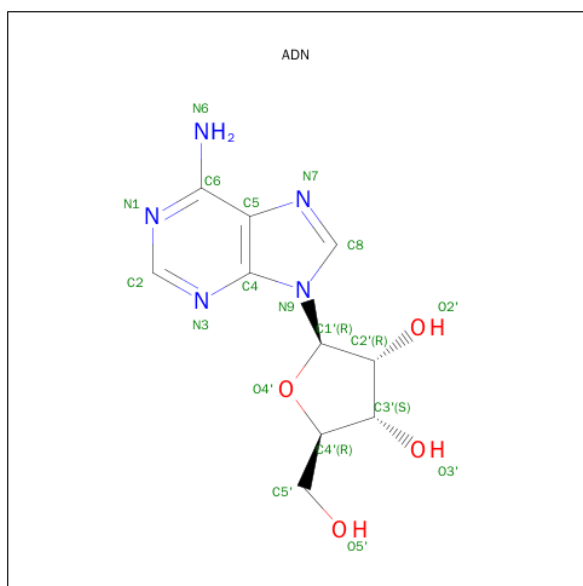
There are 5 unique types of molecules in this entry. The entry contains 9361 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ACSD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	574	Total	C	N	O	S	0	0	0
			4601	2926	834	821	20			
1	B	577	Total	C	N	O	S	0	0	0
			4626	2940	841	825	20			

- Molecule 2 is ADENOSINE (three-letter code: ADN) (formula: $C_{10}H_{13}N_5O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			19	10	5	4		
2	B	1	Total	C	N	O	0	0
			19	10	5	4		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			13	6	7		

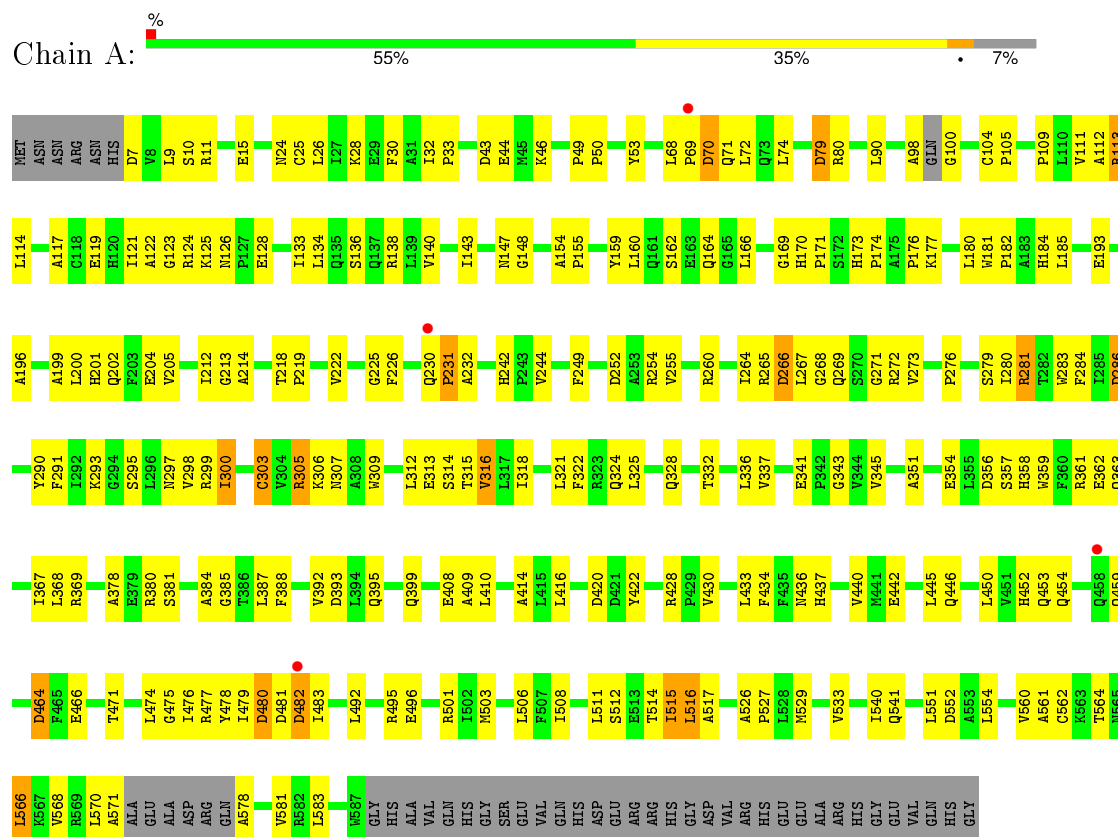
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	33	Total 33	O 33	0	0
5	B	40	Total 40	O 40	0	0

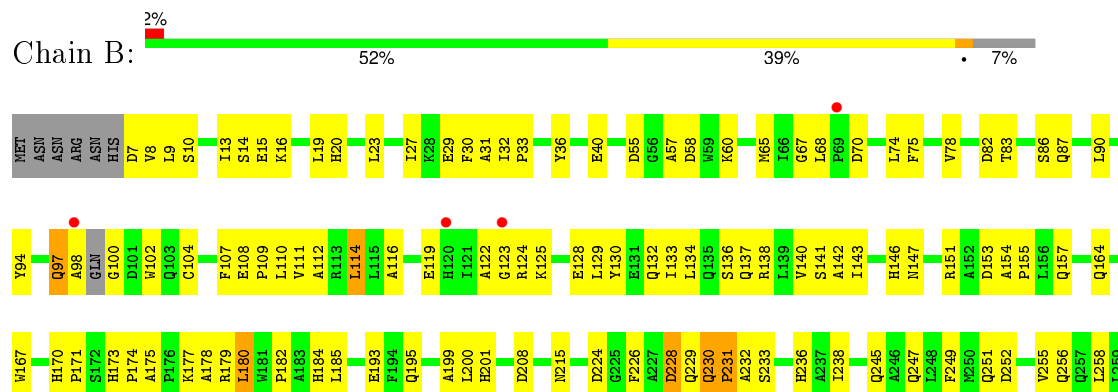
3 Residue-property plots

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: ACSD



• Molecule 1: ACSD



ARG	ARG	R436	A352	R260
ARG	L525	H437	G353	I264
HIS	A526	G438	E354	R265
GLY	P527	V439	L355	D266
ASP	L528	E442	D356	L267
VAL	Q534	P443	S357	G268
ARG	Q535	H444	F360	Q269
HIS	Q536	L445	R361	
GLU	Q541	Q446	E362	R272
GLU	Q541		V273	
ALA		L450	I367	R281
ARG	L544	Y451	L368	
HIS		H452	R369	
GLY	P547	Q459	R374	I285
VAL	L551	R463	R375	D286
GLN	D552	D464	T376	D287
HIS	I555	F465	R380	K293
GLY	A556	D472	S381	G294
	A561	D473	I382	S295
C562	L474	L474	N383	L296
K563	G475	G475	A384	N297
T564	I476	G385	G385	V298
N565	R477	I476	T386	R299
L566	Y478	F388	L387	I300
K567	I479		F388	T301
V568	D480		L394	N302
R569	D481			C303
L570	D482		Q399	V304
A571	I483		F400	R305
A572	Q489		L402	K306
GLU	S490		Y406	R307
ALA	L491		G407	A308
ASP	L492		E408	S314
R576	Y493		A409	R320
Q577	S494		L410	L321
V581	E496		D411	Q324
R582	N500		D412	D327
L583	R501		L416	Q328
P584	I502		D420	H329
W587	M503		D420	A330
GLY	L506		Q423	D331
HIS	F507		T424	T332
ALA	I508		R425	L333
VAL	L511		R428	V337
GLN	S512		P429	A340
HIS	E513		Y430	E341
GLY	T514		I431	V345
SER	I515		S432	S346
GLU	L516		F434	
VAL				A350
GLN				A351
HIS				
ASP				
GLU				

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.65Å 69.14Å 94.19Å 95.43° 101.45° 95.06°	Depositor
Resolution (Å)	91.67 – 2.95 45.95 – 2.95	Depositor EDS
% Data completeness (in resolution range)	98.2 (91.67-2.95) 94.6 (45.95-2.95)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.96Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.202 , 0.267 0.212 , 0.275	Depositor DCC
R_{free} test set	1474 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	40.2	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 29165 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	9361	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ADN, SO4, CIT

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	0.73	0/4721	0.86	6/6415 (0.1%)
1	B	0.75	0/4746	0.85	4/6448 (0.1%)
All	All	0.74	0/9467	0.86	10/12863 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	2	3
1	B	0	3
All	All	2	6

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	351	ALA	N-CA-C	7.38	130.92	111.00
1	A	474	LEU	CA-CB-CG	6.88	131.12	115.30
1	B	114	LEU	CA-CB-CG	-6.61	100.10	115.30
1	B	200	LEU	CB-CG-CD1	-6.50	99.95	111.00
1	A	566	LEU	CA-CB-CG	5.57	128.12	115.30
1	A	482	ASP	N-CA-C	5.36	125.48	111.00
1	A	303	CYS	CA-CB-SG	-5.20	104.64	114.00
1	A	266	ASP	CB-CG-OD1	5.11	122.90	118.30
1	B	394	LEU	CB-CG-CD1	-5.02	102.47	111.00
1	B	387	LEU	CB-CG-CD2	-5.01	102.48	111.00

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	351	ALA	CA
1	A	482	ASP	CA

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	268	GLY	Peptide
1	A	300	ILE	Peptide
1	A	480	ASP	Peptide
1	B	268	GLY	Peptide
1	B	327	ASP	Peptide
1	B	97	GLN	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	4601	0	4519	169	0
1	B	4626	0	4545	213	0
2	A	19	0	13	0	0
2	B	19	0	13	4	0
3	A	5	0	0	0	0
3	B	5	0	0	2	0
4	B	13	0	5	1	0
5	A	33	0	0	0	0
5	B	40	0	0	5	0
All	All	9361	0	9095	382	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 21.

All (382) 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:124:ARG:CD	1:B:571:ALA:HA	1.67	1.23
1:B:124:ARG:HD3	1:B:571:ALA:CA	1.68	1.23
1:B:563:LYS:NZ	1:B:576:ARG:HH22	1.41	1.15

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:503:MET:HE3	1:B:561:ALA:H	1.19	1.08
1:A:303:CYS:SG	1:A:305:ARG:NH2	2.29	1.05
1:B:476:ILE:HD13	1:B:492:LEU:HD13	1.32	1.04
1:B:423:GLN:HE22	1:B:536:GLN:HG3	0.92	1.03
1:B:423:GLN:NE2	1:B:536:GLN:HG3	1.75	1.01
1:B:301:THR:HG21	2:B:1588:ADN:O2'	1.64	0.98
1:B:564:THR:O	1:B:568:VAL:HG23	1.65	0.97
1:B:563:LYS:HZ2	1:B:576:ARG:HH22	1.12	0.94
1:B:563:LYS:NZ	1:B:576:ARG:NH2	2.13	0.94
1:B:29:GLU:OE1	1:B:566:LEU:HB3	1.68	0.94
1:A:134:LEU:O	1:A:138:ARG:HG2	1.73	0.88
1:B:110:LEU:HD12	1:B:114:LEU:HD12	1.56	0.88
1:A:476:ILE:HD12	1:A:479:ILE:HD12	1.55	0.87
1:B:255:VAL:HA	1:B:258:LEU:HD12	1.58	0.86
1:B:476:ILE:HG13	1:B:479:ILE:HD12	1.58	0.85
1:B:425:ARG:HG3	1:B:425:ARG:HH21	1.41	0.85
1:B:563:LYS:HZ2	1:B:576:ARG:NH2	1.73	0.85
1:B:136:SER:O	1:B:140:VAL:HG23	1.79	0.83
1:A:303:CYS:SG	1:A:305:ARG:CZ	2.67	0.83
1:B:287:ASP:HA	1:B:374:ARG:HD2	1.61	0.83
1:B:563:LYS:HZ1	1:B:576:ARG:HH22	1.25	0.82
1:A:136:SER:O	1:A:140:VAL:HG23	1.80	0.82
1:B:361:ARG:HD2	5:B:2023:HOH:O	1.81	0.80
1:A:98:ALA:HB3	1:A:100:GLY:HA2	1.63	0.79
1:A:181:TRP:HB2	1:A:182:PRO:HD2	1.64	0.79
1:B:182:PRO:HB2	1:B:184:HIS:CE1	2.18	0.79
1:B:245:GLN:OE1	1:B:296:LEU:HA	1.84	0.77
1:B:14:SER:OG	1:B:137:GLN:NE2	2.16	0.77
1:A:476:ILE:HD12	1:A:479:ILE:CD1	2.15	0.76
1:B:215:ASN:HD21	1:B:320:ARG:HB2	1.51	0.75
1:B:503:MET:HE3	1:B:561:ALA:N	1.98	0.75
1:A:299:ARG:HA	1:A:303:CYS:O	1.86	0.75
1:B:299:ARG:HA	1:B:303:CYS:O	1.86	0.74
1:A:428:ARG:HG2	1:A:540:ILE:HG12	1.68	0.74
1:B:563:LYS:HZ1	1:B:576:ARG:NH2	1.81	0.73
1:B:444:HIS:HE1	1:B:446:GLN:HB3	1.53	0.72
1:B:341:GLU:OE1	1:B:465:PHE:HB2	1.92	0.70
1:B:164:GLN:OE1	1:B:195:GLN:HA	1.91	0.70
1:B:301:THR:OG1	1:B:302:ASN:N	2.22	0.70
1:A:252:ASP:OD2	1:A:254:ARG:HB2	1.91	0.70
1:A:160:LEU:O	1:A:164:GLN:HG3	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:LEU:HD21	1:B:247:GLN:HG3	1.74	0.70
1:B:182:PRO:CB	1:B:184:HIS:CE1	2.74	0.70
1:B:387:LEU:HD21	1:B:450:LEU:HD22	1.74	0.69
1:B:388:PHE:CE1	1:B:514:THR:HG23	2.26	0.69
1:B:299:ARG:HE	1:B:569:ARG:HH22	1.38	0.69
1:B:314:SER:HB2	1:B:491:LEU:HD21	1.74	0.69
1:A:98:ALA:HB3	1:A:100:GLY:CA	2.23	0.69
1:B:167:TRP:O	1:B:177:LYS:HE3	1.93	0.69
1:A:309:TRP:HA	1:A:312:LEU:HD12	1.76	0.68
1:A:117:ALA:O	1:A:121:ILE:HD12	1.93	0.68
1:A:204:GLU:HB3	1:A:265:ARG:HB2	1.75	0.68
1:B:23:LEU:O	1:B:27:ILE:HD12	1.93	0.68
1:A:280:ILE:HG21	1:A:300:ILE:HD13	1.74	0.68
1:B:182:PRO:CB	1:B:184:HIS:HE1	2.06	0.68
1:B:445:LEU:HD12	2:B:1588:ADN:HN62	1.58	0.68
1:B:476:ILE:CG1	1:B:479:ILE:HD12	2.24	0.67
1:A:297:ASN:ND2	1:A:306:LYS:HB3	2.08	0.67
1:A:508:ILE:HD12	1:A:562:CYS:HB2	1.77	0.67
1:A:134:LEU:HB3	1:A:138:ARG:NH1	2.10	0.67
1:A:408:GLU:HG3	1:A:409:ALA:O	1.95	0.67
1:B:175:ALA:HB1	1:B:178:ALA:HB2	1.77	0.66
1:B:299:ARG:NE	1:B:569:ARG:HH22	1.94	0.66
1:A:202:GLN:NE2	1:A:271:GLY:HA2	2.11	0.66
1:A:568:VAL:HG21	1:A:578:ALA:HB1	1.78	0.65
1:B:425:ARG:HG3	1:B:425:ARG:NH2	2.09	0.65
1:A:43:ASP:OD1	1:A:44:GLU:N	2.27	0.65
1:B:382:ILE:HD12	1:B:452:HIS:CD2	2.32	0.65
1:B:9:LEU:O	1:B:13:ILE:HG13	1.97	0.65
1:A:124:ARG:NH2	1:A:571:ALA:O	2.29	0.65
1:B:108:GLU:HB3	1:B:109:PRO:HD3	1.79	0.64
1:A:297:ASN:HD21	1:A:306:LYS:HB3	1.63	0.64
1:A:541:GLN:NE2	1:A:552:ASP:OD1	2.31	0.64
1:B:444:HIS:CE1	1:B:446:GLN:HB3	2.32	0.64
1:A:225:GLY:O	1:A:272:ARG:NH2	2.26	0.64
1:A:11:ARG:NH1	1:A:15:GLU:OE2	2.31	0.64
1:A:148:GLY:HA2	1:A:392:VAL:HG11	1.79	0.64
1:A:313:GLU:CD	1:A:361:ARG:HH22	2.02	0.63
1:A:143:ILE:CD1	1:A:193:GLU:HG3	2.29	0.63
1:A:303:CYS:SG	1:A:305:ARG:NE	2.71	0.63
1:A:380:ARG:O	1:A:452:HIS:HD2	1.80	0.63
1:B:476:ILE:CD1	1:B:479:ILE:HD12	2.29	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:SER:HB3	1:A:306:LYS:HA	1.79	0.63
1:B:424:THR:HG22	1:B:428:ARG:HH21	1.63	0.62
1:B:215:ASN:ND2	1:B:320:ARG:HB2	2.13	0.62
1:A:170:HIS:HB3	1:A:173:HIS:HB2	1.81	0.62
1:B:110:LEU:CD1	1:B:114:LEU:HD12	2.29	0.61
1:B:356:ASP:O	1:B:357:SER:C	2.39	0.61
1:B:86:SER:OG	1:B:512:SER:HB2	2.00	0.61
1:A:464:ASP:OD2	1:A:466:GLU:HB3	2.00	0.61
1:A:170:HIS:HB3	1:A:173:HIS:CB	2.31	0.61
1:B:281:ARG:HD3	1:B:294:GLY:O	2.01	0.61
1:B:303:CYS:H	1:B:305:ARG:HH11	1.49	0.60
1:A:280:ILE:HG21	1:A:300:ILE:CD1	2.30	0.60
1:B:423:GLN:HE22	1:B:536:GLN:CG	1.88	0.60
1:B:357:SER:O	1:B:361:ARG:HG3	2.02	0.60
1:A:143:ILE:HD13	1:A:193:GLU:HG3	1.82	0.60
1:B:129:LEU:O	1:B:133:ILE:HG13	2.01	0.60
1:A:380:ARG:HG2	1:A:452:HIS:O	2.00	0.60
1:B:541:GLN:HA	1:B:544:LEU:HD12	1.83	0.60
1:B:476:ILE:CD1	1:B:479:ILE:CD1	2.80	0.59
1:A:180:LEU:O	1:A:298:VAL:HG13	2.01	0.59
1:B:480:ASP:OD1	1:B:480:ASP:O	2.20	0.59
1:A:430:VAL:HG21	1:A:506:LEU:HD21	1.84	0.59
1:B:424:THR:HG22	1:B:428:ARG:NH2	2.18	0.59
1:A:337:VAL:HG21	1:A:459:GLN:OE1	2.03	0.59
1:B:182:PRO:HB3	1:B:184:HIS:HE1	1.67	0.58
1:A:388:PHE:CZ	1:A:514:THR:HG23	2.38	0.58
1:B:252:ASP:HB3	1:B:255:VAL:HG23	1.85	0.58
1:A:380:ARG:O	1:A:452:HIS:CD2	2.56	0.58
1:A:68:LEU:HD12	1:A:72:LEU:CD2	2.33	0.58
1:B:424:THR:CG2	1:B:428:ARG:HH21	2.17	0.58
1:B:233:SER:O	1:B:236:HIS:HB2	2.03	0.58
1:B:308:ALA:HA	1:B:362:GLU:OE2	2.04	0.58
1:A:222:VAL:HG12	1:A:226:PHE:HE2	1.68	0.58
1:B:406:TYR:HB2	1:B:410:LEU:HD11	1.86	0.58
1:A:173:HIS:O	1:A:176:PRO:HD3	2.03	0.57
1:A:26:LEU:HD22	1:A:114:LEU:HD22	1.87	0.57
1:B:388:PHE:CZ	1:B:514:THR:HG23	2.39	0.57
1:B:94:TYR:CE1	1:B:104:CYS:HB2	2.40	0.57
1:B:98:ALA:HA	1:B:100:GLY:N	2.19	0.57
1:A:249:PHE:O	1:A:255:VAL:HG21	2.04	0.56
1:A:332:THR:HB	1:A:436:ASN:HD22	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:29:GLU:HG2	1:B:564:THR:HB	1.87	0.56
1:A:481:ASP:O	1:A:483:ILE:N	2.36	0.56
1:A:181:TRP:HB2	1:A:182:PRO:CD	2.36	0.56
1:B:157:GLN:HA	1:B:157:GLN:OE1	2.04	0.56
1:B:30:PHE:C	1:B:33:PRO:HD2	2.25	0.56
1:B:151:ARG:HG2	1:B:153:ASP:OD1	2.05	0.56
1:A:503:MET:HG2	1:A:560:VAL:HG13	1.88	0.56
1:B:110:LEU:HD12	1:B:114:LEU:CD1	2.33	0.56
1:A:9:LEU:HD11	1:B:58:ASP:HA	1.87	0.55
1:B:293:LYS:HE3	3:B:1589:SO4:O2	2.07	0.55
1:A:503:MET:CG	1:A:560:VAL:HG13	2.36	0.55
1:A:200:LEU:HD21	1:A:283:TRP:CD2	2.41	0.55
1:B:60:LYS:O	1:B:90:LEU:HD12	2.07	0.55
1:B:140:VAL:HG22	1:B:177:LYS:HG3	1.88	0.55
1:A:381:SER:HA	1:A:450:LEU:O	2.07	0.55
1:B:433:LEU:HD22	1:B:439:VAL:HB	1.89	0.55
1:B:324:GLN:O	1:B:328:GLN:HB3	2.07	0.54
1:B:481:ASP:O	1:B:483:ILE:N	2.38	0.54
1:A:356:ASP:O	1:A:357:SER:C	2.44	0.54
1:A:354:GLU:HA	1:A:357:SER:HB3	1.89	0.54
1:B:182:PRO:HB3	1:B:184:HIS:CE1	2.43	0.54
1:A:314:SER:O	1:A:318:ILE:HG12	2.08	0.54
1:A:166:LEU:O	1:A:177:LYS:NZ	2.30	0.54
1:A:169:GLY:O	1:A:171:PRO:HD3	2.08	0.54
1:B:301:THR:CG2	2:B:1588:ADN:O2'	2.47	0.54
1:B:385:GLY:HA2	1:B:445:LEU:HB3	1.89	0.54
1:B:266:ASP:OD1	1:B:267:LEU:O	2.25	0.54
1:B:201:HIS:HA	1:B:269:GLN:HA	1.90	0.54
1:A:79:ASP:HB3	1:A:90:LEU:HD11	1.89	0.54
1:B:7:ASP:HA	1:B:10:SER:OG	2.07	0.54
1:B:143:ILE:HG23	1:B:193:GLU:HB3	1.90	0.53
1:B:75:PHE:HB2	1:B:102:TRP:CH2	2.43	0.53
1:A:568:VAL:HG21	1:A:578:ALA:CB	2.38	0.53
1:B:249:PHE:CE1	1:B:255:VAL:HG11	2.44	0.53
1:B:98:ALA:HA	1:B:100:GLY:CA	2.37	0.53
1:B:226:PHE:CZ	1:B:345:VAL:HG12	2.44	0.53
1:B:15:GLU:HG2	1:B:107:PHE:CG	2.44	0.53
1:B:224:ASP:OD1	1:B:229:GLN:HB2	2.08	0.53
1:B:387:LEU:CD2	1:B:450:LEU:HD22	2.38	0.53
1:A:226:PHE:CZ	1:A:368:LEU:HG	2.44	0.53
1:B:341:GLU:HB2	1:B:465:PHE:CD2	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:LEU:CD1	1:A:478:TYR:HB3	2.39	0.53
1:A:336:LEU:HD22	1:A:433:LEU:HD11	1.90	0.53
1:A:309:TRP:O	1:A:312:LEU:HB2	2.09	0.53
1:B:299:ARG:NE	1:B:569:ARG:NH2	2.57	0.53
1:B:340:ALA:O	1:B:369:ARG:HG3	2.09	0.53
1:B:19:LEU:HD13	1:B:19:LEU:C	2.30	0.52
1:B:506:LEU:O	1:B:511:LEU:HD12	2.08	0.52
1:A:279:SER:HB3	1:A:446:GLN:NE2	2.24	0.52
1:B:173:HIS:ND1	1:B:174:PRO:HD2	2.24	0.52
1:A:385:GLY:HA2	1:A:445:LEU:HB3	1.91	0.52
1:B:376:THR:O	1:B:380:ARG:NH1	2.43	0.52
1:A:25:CYS:SG	1:A:174:PRO:HD3	2.50	0.52
1:B:124:ARG:NH1	1:B:571:ALA:O	2.43	0.52
1:A:324:GLN:O	1:A:328:GLN:HB2	2.08	0.52
1:B:55:ASP:OD1	1:B:57:ALA:N	2.37	0.52
1:A:212:ILE:HG21	1:A:219:PRO:HA	1.92	0.51
1:B:526:ALA:HB3	1:B:527:PRO:CD	2.40	0.51
1:A:266:ASP:OD1	1:A:267:LEU:O	2.29	0.51
1:B:110:LEU:CD1	1:B:114:LEU:CD1	2.89	0.51
1:B:55:ASP:C	1:B:55:ASP:OD1	2.48	0.51
1:B:446:GLN:HG3	2:B:1588:ADN:O4'	2.11	0.51
1:A:290:TYR:HB2	1:A:368:LEU:HD22	1.92	0.51
1:B:551:LEU:O	1:B:555:ILE:HG13	2.10	0.51
1:A:119:GLU:HG2	1:A:125:LYS:HA	1.91	0.51
1:B:19:LEU:HD13	1:B:19:LEU:O	2.11	0.51
1:B:30:PHE:O	1:B:33:PRO:HD2	2.11	0.51
1:B:245:GLN:OE1	1:B:297:ASN:N	2.41	0.50
1:B:508:ILE:HG21	1:B:562:CYS:SG	2.51	0.50
1:A:434:PHE:CD2	1:A:440:VAL:HG22	2.47	0.50
1:A:11:ARG:HG2	1:A:15:GLU:OE2	2.12	0.50
1:A:230:GLN:O	1:A:232:ALA:N	2.44	0.50
1:A:437:HIS:O	1:A:475:GLY:HA2	2.12	0.50
1:B:179:ARG:C	1:B:180:LEU:HG	2.31	0.50
1:B:474:LEU:O	1:B:477:ARG:HG2	2.12	0.50
1:B:124:ARG:HD3	1:B:571:ALA:HA	0.74	0.49
1:B:249:PHE:O	1:B:255:VAL:HG21	2.13	0.49
1:A:184:HIS:CE1	1:A:185:LEU:CD2	2.95	0.49
1:A:242:HIS:ND1	1:A:244:VAL:HB	2.27	0.49
1:A:170:HIS:H	1:A:176:PRO:HB3	1.78	0.49
1:A:554:LEU:HD11	1:A:560:VAL:HG22	1.93	0.49
1:A:199:ALA:HA	1:A:273:VAL:HA	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:VAL:HG12	1:A:264:ILE:CG2	2.42	0.49
1:A:387:LEU:HD21	1:A:450:LEU:HD22	1.94	0.49
1:B:576:ARG:HD2	1:B:577:GLN:H	1.76	0.49
1:A:242:HIS:CE1	1:A:244:VAL:HB	2.47	0.49
1:B:442:GLU:HG3	1:B:501:ARG:HD3	1.95	0.49
1:A:284:PHE:CE1	1:A:286:ASP:OD1	2.66	0.49
1:B:401:PHE:CD2	1:B:402:LEU:HD23	2.48	0.49
1:A:526:ALA:HB3	1:A:527:PRO:HD3	1.95	0.49
1:B:568:VAL:O	1:B:571:ALA:HB3	2.13	0.48
1:A:50:PRO:HA	1:A:53:TYR:CE1	2.48	0.48
1:B:472:ASP:OD2	1:B:494:SER:HA	2.13	0.48
1:A:202:GLN:NE2	1:A:271:GLY:CA	2.76	0.48
1:A:11:ARG:O	1:A:15:GLU:HG3	2.14	0.48
1:B:541:GLN:NE2	1:B:552:ASP:OD1	2.46	0.48
1:B:272:ARG:HD3	1:B:285:ILE:HD13	1.96	0.48
1:A:318:ILE:O	1:A:322:PHE:HD2	1.96	0.48
1:A:24:ASN:O	1:A:28:LYS:HD2	2.13	0.48
1:B:129:LEU:O	1:B:129:LEU:HD12	2.12	0.48
1:A:126:ASN:OD1	1:A:128:GLU:HB2	2.13	0.48
1:B:476:ILE:HD11	1:B:479:ILE:CD1	2.43	0.48
1:A:428:ARG:HG3	1:A:540:ILE:HD11	1.96	0.48
1:B:337:VAL:HG21	1:B:459:GLN:HG2	1.95	0.48
1:B:111:VAL:O	1:B:112:ALA:C	2.53	0.48
1:A:260:ARG:O	1:A:260:ARG:HG2	2.13	0.48
1:A:79:ASP:CB	1:A:90:LEU:HD11	2.43	0.47
1:A:564:THR:HG21	1:A:581:VAL:HG13	1.95	0.47
1:B:547:PRO:O	5:B:2037:HOH:O	2.20	0.47
1:A:503:MET:HE3	1:A:561:ALA:HB2	1.95	0.47
1:A:74:LEU:C	1:A:74:LEU:HD23	2.35	0.47
1:B:489:GLN:O	1:B:489:GLN:HG3	2.13	0.47
1:A:511:LEU:O	1:A:512:SER:C	2.51	0.47
1:A:503:MET:HG2	1:A:560:VAL:CG1	2.45	0.47
1:B:170:HIS:ND1	1:B:171:PRO:HD2	2.29	0.47
1:B:228:ASP:O	1:B:231:PRO:HD2	2.14	0.47
1:B:577:GLN:NE2	5:B:2039:HOH:O	2.46	0.47
1:B:226:PHE:CE2	1:B:368:LEU:HG	2.49	0.47
1:A:515:ILE:HG22	1:A:516:LEU:N	2.30	0.47
1:A:134:LEU:HB3	1:A:138:ARG:HH12	1.79	0.47
1:B:382:ILE:HD11	1:B:452:HIS:HE2	1.78	0.47
1:A:526:ALA:HB3	1:A:527:PRO:CD	2.44	0.47
1:A:393:ASP:OD1	1:A:395:GLN:HB2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:303:CYS:H	1:B:305:ARG:NH1	2.12	0.47
1:B:170:HIS:HB3	1:B:173:HIS:HB3	1.97	0.47
1:B:321:LEU:HD13	1:B:478:TYR:HB3	1.97	0.46
1:A:471:THR:HG22	1:A:495:ARG:HA	1.96	0.46
1:A:196:ALA:O	1:A:276:PRO:HD2	2.15	0.46
1:B:68:LEU:O	1:B:70:ASP:N	2.49	0.46
1:B:97:GLN:O	1:B:98:ALA:HB3	2.16	0.46
1:B:576:ARG:CD	1:B:577:GLN:H	2.29	0.46
1:A:281:ARG:NH1	1:A:293:LYS:HE2	2.31	0.45
1:B:258:LEU:HB3	1:B:264:ILE:HG12	1.98	0.45
1:A:422:TYR:HE2	1:A:450:LEU:HD13	1.81	0.45
1:A:122:ALA:HB3	1:A:570:LEU:HD23	1.98	0.45
1:A:154:ALA:N	1:A:155:PRO:HD2	2.32	0.45
1:B:256:GLN:HB3	1:B:260:ARG:HH22	1.82	0.45
1:B:208:ASP:HB3	1:B:350:ALA:HB2	1.98	0.45
1:A:214:ALA:HB2	1:A:222:VAL:HG21	1.99	0.45
1:B:116:ALA:HA	1:B:119:GLU:OE1	2.15	0.45
1:A:184:HIS:CE1	1:A:185:LEU:HD23	2.51	0.45
1:A:143:ILE:O	1:A:147:ASN:ND2	2.42	0.45
1:B:32:ILE:N	1:B:33:PRO:CD	2.80	0.45
1:B:416:LEU:HD11	1:B:528:LEU:HD13	1.97	0.45
1:B:382:ILE:HD12	1:B:452:HIS:HD2	1.81	0.45
1:A:25:CYS:HB3	1:A:566:LEU:HD23	1.98	0.45
1:B:430:VAL:HG21	1:B:443:PRO:HG3	1.98	0.45
1:B:15:GLU:O	1:B:19:LEU:HB2	2.15	0.44
1:A:111:VAL:HG11	1:A:133:ILE:HG21	1.99	0.44
1:B:170:HIS:CG	1:B:173:HIS:HB2	2.52	0.44
1:A:159:TYR:O	1:A:162:SER:OG	2.32	0.44
1:B:306:LYS:HD3	1:B:362:GLU:HG2	2.00	0.44
1:A:321:LEU:O	1:A:325:LEU:HD12	2.17	0.44
1:A:49:PRO:O	1:A:50:PRO:C	2.56	0.44
1:B:576:ARG:CD	1:B:577:GLN:N	2.80	0.44
1:B:525:LEU:HD23	1:B:525:LEU:HA	1.59	0.44
1:A:416:LEU:O	1:A:420:ASP:HB2	2.17	0.44
1:A:30:PHE:HD1	1:A:121:ILE:HG21	1.83	0.44
1:B:134:LEU:O	1:B:138:ARG:HG3	2.16	0.44
1:A:529:MET:O	1:A:533:VAL:HG23	2.18	0.44
1:B:576:ARG:CG	1:B:577:GLN:N	2.80	0.44
1:B:107:PHE:CE2	1:B:137:GLN:HG3	2.52	0.44
1:B:23:LEU:HD23	1:B:23:LEU:HA	1.82	0.44
1:A:199:ALA:HA	1:A:272:ARG:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:GLN:HB3	1:A:231:PRO:HD3	1.99	0.44
1:B:416:LEU:O	1:B:420:ASP:HB2	2.17	0.44
1:B:496:GLU:O	1:B:500:ASN:OD1	2.35	0.44
1:B:511:LEU:O	1:B:512:SER:C	2.56	0.44
1:B:153:ASP:N	1:B:153:ASP:OD1	2.50	0.43
1:A:307:ASN:O	1:A:362:GLU:HB2	2.18	0.43
1:A:7:ASP:HA	1:A:10:SER:HB2	1.99	0.43
1:A:399:GLN:NE2	1:A:517:ALA:O	2.52	0.43
1:B:360:PHE:C	1:B:362:GLU:N	2.71	0.43
1:B:412:ASP:OD1	1:B:522:ARG:NH1	2.52	0.43
1:A:79:ASP:O	1:A:80:ARG:HB3	2.18	0.43
1:B:346:SER:HB2	1:B:361:ARG:HA	2.00	0.43
1:A:293:LYS:HB3	1:A:367:ILE:HB	2.00	0.43
1:A:70:ASP:O	1:A:71:GLN:HB2	2.17	0.43
1:B:299:ARG:C	1:B:300:ILE:HG12	2.39	0.43
1:B:401:PHE:HD2	1:B:402:LEU:HD23	1.84	0.43
1:A:453:GLN:O	1:A:454:GLN:HG3	2.18	0.43
1:B:36:TYR:HB3	1:B:67:GLY:O	2.18	0.43
1:B:74:LEU:C	1:B:74:LEU:HD23	2.39	0.43
1:A:315:THR:HG23	1:A:341:GLU:OE1	2.19	0.43
1:B:503:MET:HE2	1:B:561:ALA:CB	2.48	0.43
1:B:331:ASP:N	1:B:331:ASP:OD1	2.51	0.43
1:B:329:HIS:O	1:B:333:LEU:HG	2.19	0.43
1:A:122:ALA:CB	1:A:570:LEU:HD23	2.49	0.43
1:A:111:VAL:O	1:A:112:ALA:C	2.57	0.43
1:A:422:TYR:CE2	1:A:450:LEU:HD13	2.54	0.42
1:A:345:VAL:O	1:A:345:VAL:HG13	2.18	0.42
1:A:309:TRP:CE2	1:A:358:HIS:CE1	3.07	0.42
1:A:410:LEU:HB3	1:A:414:ALA:HB3	2.01	0.42
1:B:245:GLN:OE1	1:B:296:LEU:CA	2.62	0.42
1:B:82:ASP:OD2	1:B:86:SER:HB2	2.19	0.42
1:A:385:GLY:CA	1:A:445:LEU:HB3	2.49	0.42
1:B:380:ARG:HG2	5:B:2029:HOH:O	2.19	0.42
1:B:40:GLU:O	1:B:65:MET:HG3	2.19	0.42
1:B:83:THR:HB	1:B:584:PRO:O	2.19	0.42
1:A:476:ILE:CD1	1:A:479:ILE:CD1	2.93	0.42
1:B:541:GLN:HA	1:B:544:LEU:CD1	2.47	0.42
1:A:291:PHE:HB2	1:A:369:ARG:HB3	2.02	0.42
1:A:32:ILE:N	1:A:33:PRO:CD	2.83	0.42
1:B:383:MET:HG2	1:B:384:ALA:N	2.33	0.42
1:B:438:GLY:HA3	1:B:475:GLY:H	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:381:SER:HA	1:B:450:LEU:O	2.20	0.42
1:A:218:THR:O	1:A:219:PRO:C	2.56	0.42
4:B:1590:CIT:H41	4:B:1590:CIT:O1	2.19	0.42
1:A:313:GLU:O	1:A:316:VAL:HB	2.19	0.42
1:B:128:GLU:HG3	1:B:132:GLN:NE2	2.35	0.42
1:B:146:HIS:HD2	1:B:147:ASN:ND2	2.18	0.42
1:B:130:TYR:CE2	1:B:134:LEU:HD11	2.54	0.42
1:B:564:THR:CG2	1:B:581:VAL:HG22	2.50	0.41
1:A:299:ARG:HB2	1:A:299:ARG:NH1	2.35	0.41
1:B:437:HIS:O	1:B:475:GLY:HA2	2.19	0.41
1:A:359:TRP:O	1:A:363:GLN:HG2	2.20	0.41
1:A:213:GLY:O	1:A:343:GLY:HA2	2.20	0.41
1:B:360:PHE:O	1:B:362:GLU:N	2.53	0.41
1:B:119:GLU:HG2	1:B:125:LYS:HA	2.02	0.41
1:B:16:LYS:HE3	1:B:20:HIS:CE1	2.55	0.41
1:B:503:MET:CE	1:B:561:ALA:CB	2.98	0.41
1:B:129:LEU:HD22	1:B:569:ARG:HB2	2.01	0.41
1:B:428:ARG:O	1:B:432:SER:OG	2.25	0.41
1:A:201:HIS:HA	1:A:269:GLN:HA	2.03	0.41
1:B:251:GLN:NE2	5:B:2015:HOH:O	2.38	0.41
1:A:318:ILE:HG23	1:A:322:PHE:HE2	1.86	0.41
1:A:318:ILE:CG2	1:A:322:PHE:CE2	3.04	0.41
1:B:463:ARG:O	1:B:464:ASP:CB	2.69	0.41
1:A:508:ILE:HD13	1:A:583:LEU:HD12	2.03	0.41
1:A:476:ILE:HD13	1:A:492:LEU:HD13	2.01	0.41
1:B:293:LYS:HB3	1:B:367:ILE:HB	2.01	0.41
1:B:433:LEU:HD22	1:B:439:VAL:CB	2.50	0.41
1:A:104:CYS:HA	1:A:105:PRO:HD2	1.90	0.41
1:A:109:PRO:O	1:A:113:ARG:HB2	2.20	0.41
1:B:230:GLN:O	1:B:232:ALA:N	2.54	0.41
1:A:69:PRO:HD2	1:A:121:ILE:HD11	2.02	0.41
1:B:27:ILE:HA	1:B:31:ALA:HB3	2.03	0.41
1:A:44:GLU:HG3	1:A:46:LYS:H	1.86	0.41
1:B:442:GLU:HG3	1:B:501:ARG:HH11	1.86	0.41
1:B:199:ALA:HA	1:B:273:VAL:HA	2.02	0.41
1:B:122:ALA:C	1:B:124:ARG:H	2.24	0.40
1:A:182:PRO:HB3	1:A:184:HIS:ND1	2.36	0.40
1:A:541:GLN:HB2	1:A:551:LEU:CD2	2.51	0.40
1:B:534:GLN:NE2	1:B:555:ILE:O	2.54	0.40
1:B:154:ALA:N	1:B:155:PRO:CD	2.84	0.40
1:B:435:PHE:HB2	1:B:544:LEU:HD22	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:ALA:O	1:A:387:LEU:N	2.54	0.40
1:A:321:LEU:HD11	1:A:478:TYR:CB	2.51	0.40
1:A:267:LEU:H	1:A:267:LEU:HG	1.61	0.40
1:B:534:GLN:HG3	1:B:587:TRP:CE2	2.56	0.40
1:B:232:ALA:HB2	1:B:238:ILE:HG13	2.03	0.40
1:B:369:ARG:NH2	3:B:1589:SO4:O2	2.53	0.40
1:A:321:LEU:HD11	1:A:478:TYR:HB3	2.04	0.40
1:A:133:ILE:HG12	1:A:174:PRO:O	2.22	0.40
1:A:442:GLU:HG3	1:A:501:ARG:HH11	1.86	0.40
1:B:78:VAL:HG11	1:B:87:GLN:HG3	2.02	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	568/620 (92%)	501 (88%)	59 (10%)	8 (1%)	14	49
1	B	571/620 (92%)	494 (86%)	65 (11%)	12 (2%)	9	37
All	All	1139/1240 (92%)	995 (87%)	124 (11%)	20 (2%)	11	42

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	482	ASP
1	B	351	ALA
1	B	482	ASP
1	A	464	ASP
1	A	515	ILE
1	A	516	LEU
1	B	141	SER
1	B	464	ASP

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Mol	Chain	Res	Type
1	B	516	LEU
1	A	378	ALA
1	B	142	ALA
1	B	228	ASP
1	B	399	GLN
1	B	556	ALA
1	A	316	VAL
1	A	231	PRO
1	B	123	GLY
1	B	231	PRO
1	B	515	ILE
1	A	123	GLY

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	485/522 (93%)	476 (98%)	9 (2%)	65	89
1	B	487/522 (93%)	474 (97%)	13 (3%)	52	84
All	All	972/1044 (93%)	950 (98%)	22 (2%)	58	86

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	ASP
1	A	79	ASP
1	A	113	ARG
1	A	281	ARG
1	A	286	ASP
1	A	305	ARG
1	A	477	ARG
1	A	480	ASP
1	A	496	GLU
1	B	8	VAL
1	B	180	LEU

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Mol	Chain	Res	Type
1	B	230	GLN
1	B	272	ARG
1	B	281	ARG
1	B	362	GLU
1	B	375	ARG
1	B	481	ASP
1	B	541	GLN
1	B	569	ARG
1	B	570	LEU
1	B	576	ARG
1	B	582	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	20	HIS
1	B	137	GLN
1	B	324	GLN
1	B	423	GLN
1	B	500	ASN
1	B	577	GLN

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](#)

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	ADN	A	1588	-	16,21,21	1.22	1 (6%)	16,31,31	2.43	2 (12%)
3	SO4	A	1589	-	4,4,4	0.14	0	6,6,6	0.64	0
2	ADN	B	1588	-	16,21,21	1.12	2 (12%)	16,31,31	2.33	1 (6%)
3	SO4	B	1589	-	4,4,4	0.29	0	6,6,6	0.51	0
4	CIT	B	1590	-	3,12,12	1.44	1 (33%)	3,17,17	1.63	1 (33%)

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	ADN	A	1588	-	-	0/2/22/22	0/3/3/3
3	SO4	A	1589	-	-	0/0/0/0	0/0/0/0
2	ADN	B	1588	-	-	0/2/22/22	0/3/3/3
3	SO4	B	1589	-	-	0/0/0/0	0/0/0/0
4	CIT	B	1590	-	-	0/6/16/16	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1588	ADN	C5-N7	-2.03	1.32	1.39
4	B	1590	CIT	O7-C3	2.43	1.47	1.43
2	A	1588	ADN	O4'-C1'	3.20	1.45	1.41
2	B	1588	ADN	O4'-C1'	3.32	1.45	1.41

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1588	ADN	N3-C2-N1	-9.01	122.00	128.89
2	B	1588	ADN	N3-C2-N1	-8.53	122.36	128.89
2	A	1588	ADN	C4-C5-N7	-2.18	107.48	109.48
4	B	1590	CIT	C3-C2-C1	2.01	118.18	114.96

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1588	ADN	4	0
3	B	1589	SO4	2	0
4	B	1590	CIT	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 ⓘ

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	574/620 (92%)	-0.12	4 (0%) 89 76	8, 20, 31, 38	0
1	B	577/620 (93%)	-0.04	12 (2%) 67 46	8, 20, 31, 46	2 (0%)
All	All	1151/1240 (92%)	-0.08	16 (1%) 78 59	8, 20, 31, 46	2 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	482	ASP	4.8
1	B	571	ALA	4.4
1	B	69	PRO	4.2
1	B	409	ALA	3.4
1	B	352	ALA	3.0
1	B	407	GLY	3.0
1	A	69	PRO	2.8
1	B	123	GLY	2.8
1	B	98	ALA	2.6
1	B	351	ALA	2.5
1	B	354	GLU	2.4
1	A	458	GLN	2.4
1	B	350	ALA	2.3
1	B	120	HIS	2.1
1	A	230	GLN	2.1
1	B	481	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [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
2	ADN	A	1588	19/19	0.92	0.22	1.36	42,43,45,46	0
2	ADN	B	1588	19/19	0.95	0.19	0.23	30,31,36,38	0
3	SO4	A	1589	5/5	0.97	0.15	-0.29	19,24,25,25	0
4	CIT	B	1590	13/13	0.95	0.15	-0.95	16,29,32,33	0
3	SO4	B	1589	5/5	0.99	0.11	-3.08	18,21,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.